

# ParaView User's Guide (v3.10)

How to unleash the beast!

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# Introduction

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## About Paraview

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### What is ParaView?

ParaView is an open-source, multi-platform application for the visualization and analysis of scientific data sets, primarily those that are defined natively in a two- or three-dimensional space including those that extend into the temporal dimension.

The front end graphical user interface (GUI) has an open, flexible and intuitive user interface that still gives you fine grained and open ended control of the data manipulation and display processing needed to explore and present complex data as you see fit.

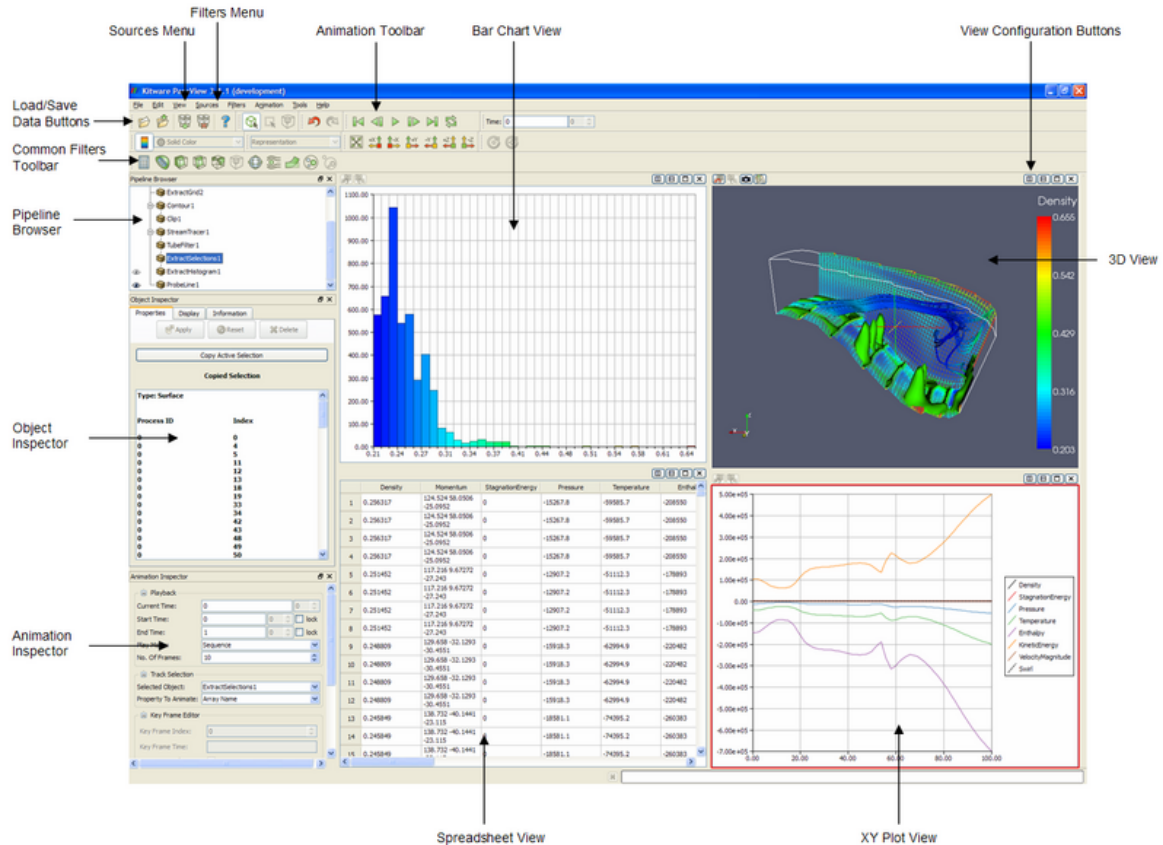
ParaView has extensive scripting and batch processing capabilities. The standard scripting interface uses the widely used python programming language for scripted control. As with the GUI, the python scripted control is easy to learn, notably including the ability to record actions in the GUI and save them out as succinct human readable python programs and powerful, including the ability to write scripted filters that run on the server that have access to every bit of your data on a large parallel machine.

ParaView's data processing and rendering components are built upon a modular and scalable distributed-memory parallel architecture in which many processors operate synchronously on different portions of the data. ParaView's scalable architecture allows you to run ParaView directly on anything from a small netbook class machine up to the world largest supercomputer. The size of the data sets ParaView can handle in practice though varies widely depending on the size of the machine that ParaView's server components are run. Thus people frequently do both, taking advantage of ParaView's client/server architecture to connect to and control the supercomputer from the netbook.

ParaView is meant to be easily extended and customized into new applications and be used by or make use of other tools. Correspondingly there are a number of different interfaces to ParaView's data processing and visualization engine, for example ParaViewWeb <sup>[1]</sup> a web based one. This book will not cover any of those or describe in great detail the programmers interface to the ParaView engine. We will instead focus our attention on understanding the standard ParaView GUI based application.

## User Interface

The different sections of ParaView's Graphical User Interface (GUI) are shown below. Of particular importance in the following discussion are the File and Filter menus which allow one to open files and manipulate data, the Pipeline Browser which displays the Visualization Pipeline, the Object Inspector with its Properties, Display and Information tabs where one can control any given module within the pipeline, and the View area where data is displayed in one or more windows.



## Modality

One very important thing to keep in mind when using ParaView is that the GUI is very modal. At any given time you will have one "active" module within the Visualization Pipeline, one "active" View, and one "active" Selection. For example when you click on the name of a reader or source within the Pipeline Browser, it becomes the active module and the properties of that filter are displayed in the Object Inspector. Likewise when you click within a different View, that view becomes the Active View and the visibility "eye" icons in the Pipeline Browser are changed to show what filters are displayed within this View. These concepts will be described in detail in later chapters [2][3][4]. For now you should be aware that the information displayed in the GUI always pertains to these active entities.

## Basics of Visualization

Put simply, the process of visualization is taking raw data and converting it to a form that is viewable and understandable to humans. This allows us to get a better cognitive understanding of our data. Scientific visualization is specifically concerned with the type of data that has a well-defined representation in 2D or 3D space. Data that comes from simulation meshes and scanner data is well suited for this type of analysis.

There are three basic steps to visualizing your data: reading, filtering, and rendering. First, your data must be read into ParaView. Next, you may apply any number of filters that process the data to generate, extract, or derive features from the data. Finally, a viewable image is rendered from the data and you can then change the viewing parameters or rendering modality for best visual effect.

## The Pipeline Concept

In ParaView, these steps are made manifest in a Visualization Pipeline. That is one visualizes data by building up a set of modules, each of which takes in some data, operates on it, and presents the result as a new dataset. This begins with a Reader module whose task it is to ingest data off of files on disk.

Reading data into ParaView is often as simple as selecting Open from the File menu, and then clicking the glowing Accept button on the reader's Object Inspector tab. ParaView comes with support for a large number of file formats[5], and its modular architecture makes it possible to add new file readers[6].

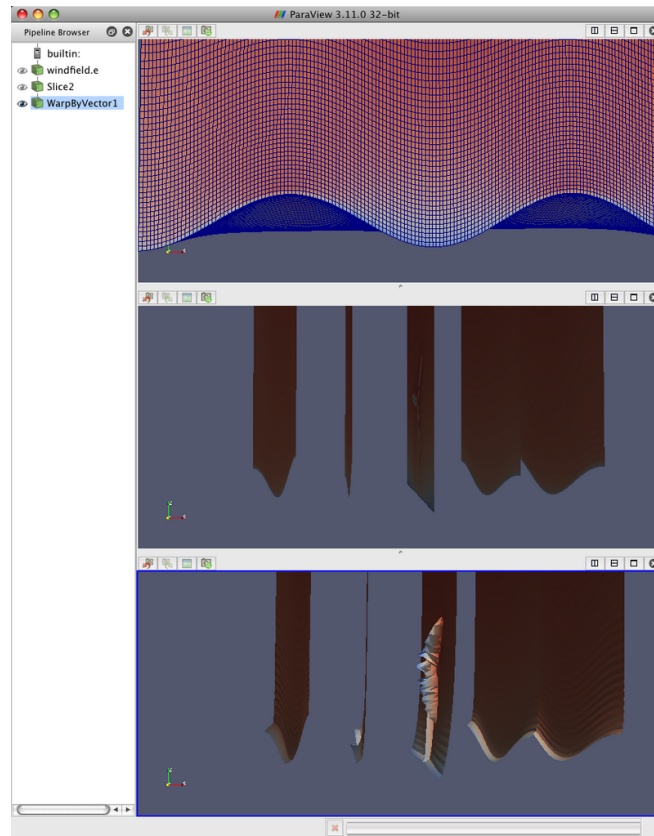
Once a file is read, ParaView automatically renders it in a View. In ParaView, a View is simply a window that shows data. There are different types of Views, ranging from qualitative computer graphics rendering of the data to quantitative spreadsheet presentations of the data values as text. ParaView picks a suitable view type for your data automatically, but you are free to change the view type, modify the rendering parameters of the data in the view, and even create new views simultaneously as you see fit to better understand what you've read in. Additionally, high level meta information about the data including names, types and ranges of arrays, temporal ranges, memory size and geometric extent can be found in the Information Tab.

One can learn a great deal about a given dataset with a one element Visualization Pipeline consisting of just a reader module. You can learn much more, and in fact perform arbitrarily complex analyses and data manipulations, adding modules to build up the pipeline. In ParaView you can create arbitrarily complex visualization pipelines, including ones with multiple readers, merging and branching pipelines. You build up a pipeline by choosing the next filter in the sequence from the Filters menu. Once you click accept, this new filter will read in the data produced by the formerly active filter and perform some processing on that data. The new filter then becomes the active one. Filters then are created differently from but operate in the same manner as readers. At all points you use the Pipeline Inspector to choose the active filter and then the Object Inspector to configure it.

The Pipeline Browser is where the overall Visualization Pipeline is displayed and controllable from. The Object Inspector is where the specific parameters of one particular module within the Pipeline are displayed and controllable from. The Object inspector has three tabs, one presents the parameters of the processing done within that module, another presents the parameters of how the output of that module will be displayed in a View, the last presents the meta information about the data produced by the module as described above.

The following figure demonstrates a three element visualization pipeline, where the output of each module in the pipeline is displayed in its own view. Here a reader takes in a vector field, defined on a curvilinear grid, which comes from a simulation study of a wind turbine. Next a slice filter produces slices of the field on five equally spaced planes along the X axis. Finally a warp filter warps those planes along the direction of the vector field, which primarily moves the planes downwind but also shows some complexity at the location of the wind turbine.





There are more than one hundred filters available to choose from in total, all of which manipulate the data in different ways. The full list of filters is available in the Appendix <sup>[5]</sup> and within the application under the help menu. Note that many of the filters in the menu will be grayed out and not selectable at any given time. That is because any given filter may only operate on particular types of data. For example the Extract Subset will only operate on structured data sets so it is only enabled when the module you are building on top of produces image data, rectilinear grid data, or structured grid data. (This input restrictions are also listed in the Appendix <sup>[7]</sup> and help menu). In this situation you can often find a similar filter which does accept your data, or apply a filter which transforms you data into the required format. Since ParaView 3.10 you can ask ParaView to try to do the conversion for you automatically, by clicking "Auto convert Properties" in the application settings <sup>[8]</sup>. The mechanics of applying filters are described fully in the Manipulating Data <sup>[9]</sup> chapter.

## Making Mistakes

Frequently, new users of ParaView falter when they open their data, or apply a filter, and do not see it immediately because they have not pressed the Apply button. ParaView was designed to operate on large data sets, for which any given operation could take a long time to finish, in this situation you need the Apply button/commit behavior so that you have a chance to be confident of your change before it takes effect. The highlighted Apply button is a reminder that the parameters of one or more filters are out of sync with the data that you are viewing. Hitting the Apply button accepts your change (or changes) whereas hitting the Reset button reverts the options back to the last time you hit Apply. If you are working with small data sets, you may want to turn off this behavior with the Auto Accept setting under the Application's settings <sup>[8]</sup>.

The Apply behavior circumvents a great number of mistakes but not all of them. If you make some change to a filter or to the pipeline itself and later find that you are not satisfied with the result, hit the Undo button. You can undo all the way back to the start of your ParaView session and redo all the way forward if you like. You can also undo and redo camera motion by using the camera undo and redo buttons located above each View window.

## Persistent Sessions

If on the other hand you are satisfied with your visualization results, you may want to save your work so that you can return to it at some future time. You can do so by using ParaView's Save State (File->Save State) and Save Trace (Tools -> Save Trace) features. In either case, ParaView produces human readable text files (XML files for State and Python script for Trace) that can be restored or modified and restored later. This is very useful for batch processing, which is discussed in the Python Scripting <sup>[10]</sup> chapter.

To save state means to save enough information about the ParaView session to restore it later and thus show exactly the same result. ParaView does so by saving the current visualization pipeline and the parameters of the filters within it.

If you turn on a trace recording when you first start using ParaView, saving a trace can be used for the same purpose as saving state. However, a trace records all of your actions, including the ones that you later undo, as you do them. It is a more exact recording then of not only what you did, but how you did it. Traces are saved as python scripts, which ParaView can play back in either batch mode or within an interactive GUI session. You can use traces then to automate repetitive tasks by recording just that action. It is also an ideal tool to learn ParaView's python scripting API.

## Client/Server Visualization

With small datasets it is usually quite sufficient to run ParaView as a single process on a small laptop or desktop class machine. For large data sets, a single machine is not likely to have enough processing power and much more importantly memory to process the data. In this situation one runs an MPI parallel ParaView Server process on a large machine to do computationally and memory expensive data processing and/or rendering tasks and then connects to that server within the familiar GUI application.

When connected to a remote server the only difference you will see will be that the Visualization Pipeline displayed in the Pipeline Browser will begin with the name of the server you are connected to rather than the word 'builtin' which indicates that you are connected to a virtual server residing in the same process as the GUI. When connected to a remote server, the File Open dialog presents the list of files that live on the remote machine's file system rather than the client's. Depending on the server's capabilities and on the data size and your application settings (Edit->Settings->Render View->Server) the data will either be rendered remotely and pixels will be sent to the client or the geometry will be delivered and rendered locally. Large data visualization is described fully in the Client Server Visualization <sup>[11]</sup> Chapter.

## References

- [1] <http://www.paraview.org/Wiki/ParaViewWeb>
- [2] [http://paraview.org/Wiki/ParaView/Displaying\\_Data#Multiple\\_Views](http://paraview.org/Wiki/ParaView/Displaying_Data#Multiple_Views)
- [3] [http://paraview.org/Wiki/ParaView/UsersGuide/Filtering\\_Data#Pipeline\\_Basics](http://paraview.org/Wiki/ParaView/UsersGuide/Filtering_Data#Pipeline_Basics)
- [4] [http://paraview.org/Wiki/ParaView/Users\\_Guide/Selection](http://paraview.org/Wiki/ParaView/Users_Guide/Selection)
- [5] [http://paraview.org/Wiki/ParaViewUsersGuide/List\\_of\\_readers](http://paraview.org/Wiki/ParaViewUsersGuide/List_of_readers)
- [6] [http://paraview.org/Wiki/Writing\\_ParaView\\_Readers](http://paraview.org/Wiki/Writing_ParaView_Readers)
- [7] [http://paraview.org/Wiki/ParaViewUsersGuide/List\\_of\\_filters](http://paraview.org/Wiki/ParaViewUsersGuide/List_of_filters)
- [8] [http://paraview.org/Wiki/ParaView/Users\\_Guide/Settings](http://paraview.org/Wiki/ParaView/Users_Guide/Settings)
- [9] [http://paraview.org/Wiki/ParaView/UsersGuide/Filtering\\_Data](http://paraview.org/Wiki/ParaView/UsersGuide/Filtering_Data)
- [10] [http://paraview.org/Wiki/ParaView/Python\\_Scripting](http://paraview.org/Wiki/ParaView/Python_Scripting)
- [11] [http://paraview.org/Wiki/Users\\_Guide\\_Client-Server\\_Visualization](http://paraview.org/Wiki/Users_Guide_Client-Server_Visualization)

# Loading Data

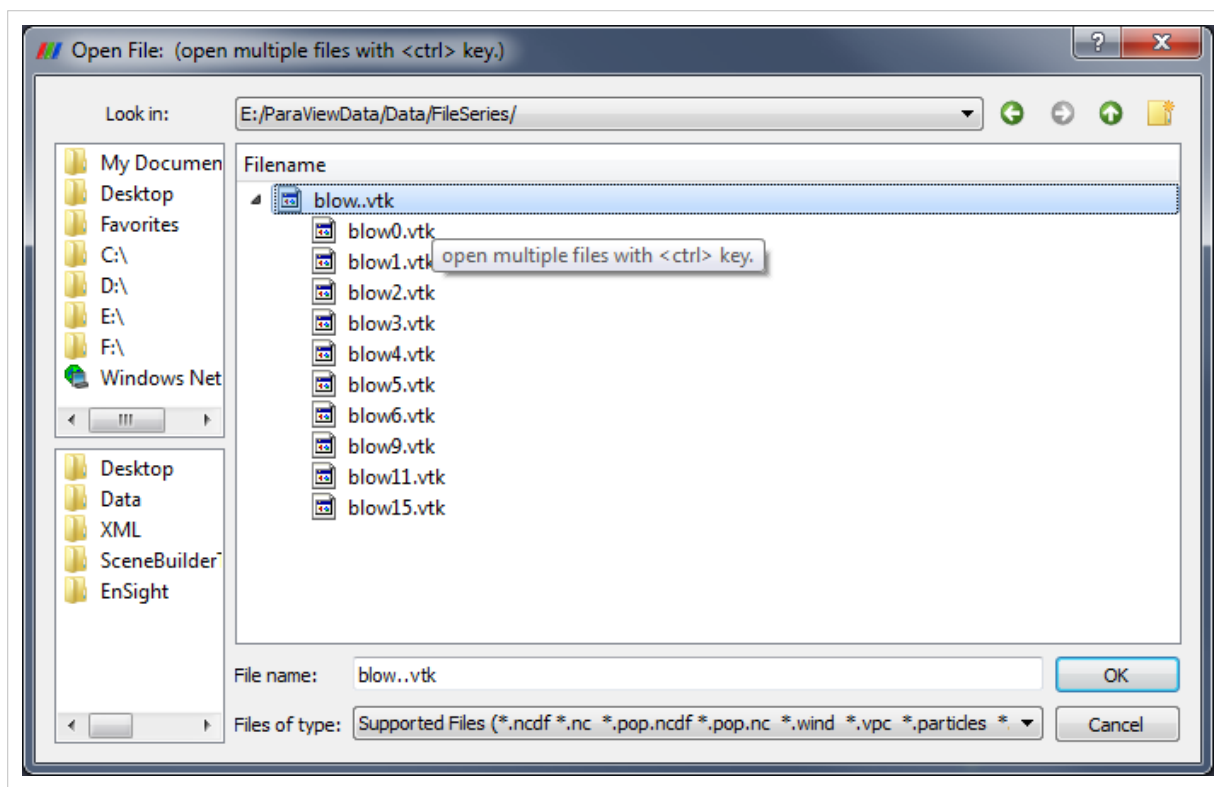
## Data Ingestion

### Introduction

Loading data is a fundamental operation in using ParaView for visualization. As you would expect, the **Open** option from the **File** menu and the **Open Button** from the toolbar both allow you to load data into ParaView. ParaView understands many scientific data file formats. The most comprehensive list is given in the List of Readers <sup>[5]</sup> appendix. Because of ParaView's modular design it is easy to integrate new readers. If the formats you care about are not listed there ask the mailing list first to see if anyone has a reader for the format or, if you want to create your own readers for ParaView see the Plugin HowTo <sup>[1]</sup> section and the Writing Readers <sup>[6]</sup> appendix of this book.

### Opening File / Time Series

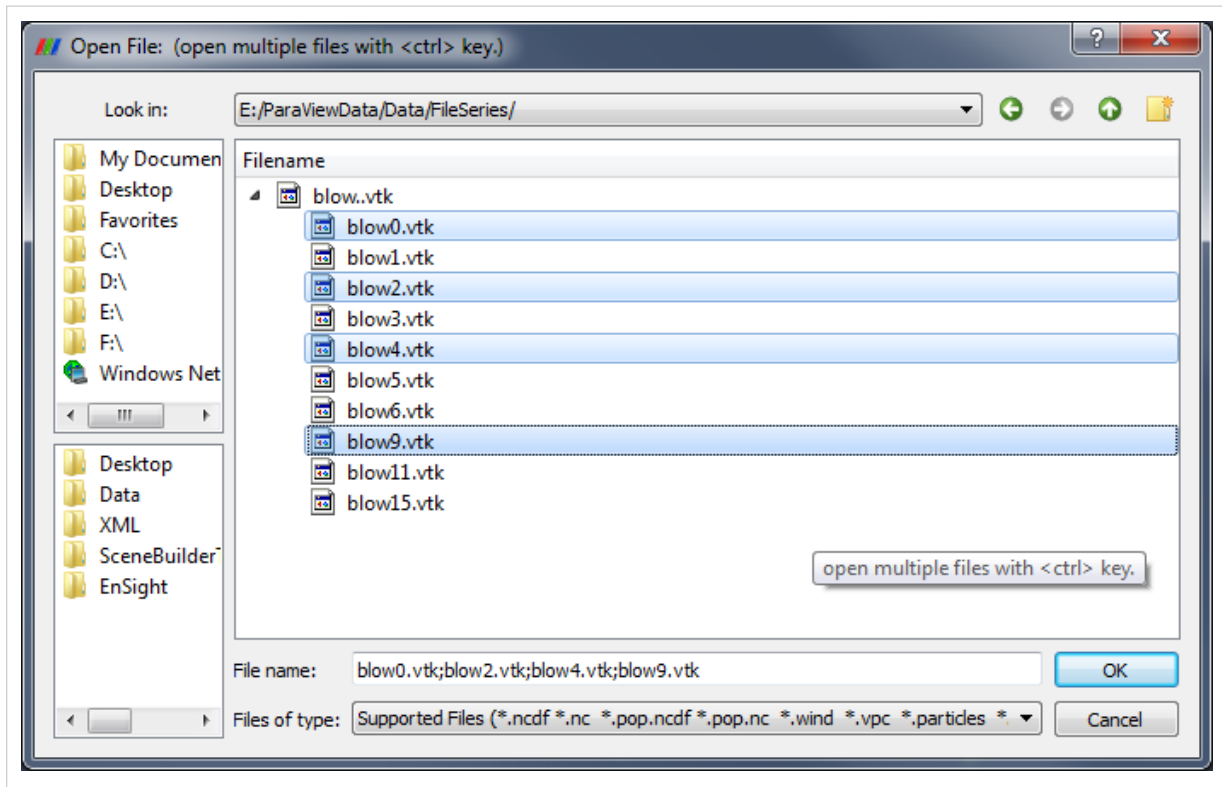
ParaView recognizes file series by using certain patterns in the name of files including: fooN.vtk foo\_N.vtk foo-N.vtk foo.N.vtk Nfoo.vtk N.foo.vtk foo.vtk.N foo.vtk-sN Where N is an integer (with any number of leading zeros). To load a file series, first make sure that the file names match one of the patterns described above. Next, navigate to the directory where the file series is. The file browser should look like this:



You can expand the file series by clicking on the triangle, as shown in the above diagram. Simply select the group (in the picture named blow..vtk) and click ok. The reader will store all the filenames and treat each file as a time step. You can now animate, use annotate time filter, do anything you can do with readers that natively support time. If you want to load a single step of a file series just expand the triangle and select the file you are interested in.

## Opening Multiple Files

ParaView supports loading multiple files as long as they exist in the same directory. Just hold the ctrl key down while selecting each file, or hold shift to select all files in a range.



## State Files

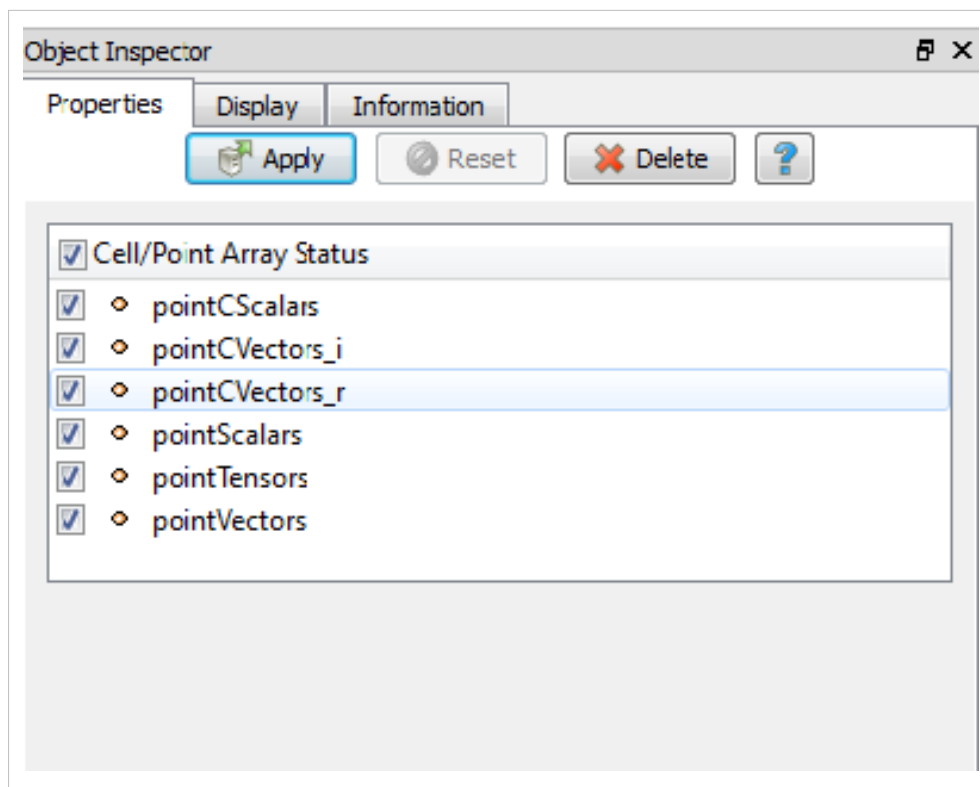
Another option is to load a previously saved state file (**File** menu, **Load State**). This will return ParaView to its state at the time the file was saved by loading data files, applying filters.

## Advanced Data Loading

If you commonly load the same data into ParaView each time, you can streamline the process by launching ParaView with the data command-line argument (`--data=data_file`). You can also launch ParaView

## Object Inspector

Note that opening a file is a two step process, so that you do not see any data after opening a data file. Instead, you see that the object inspector is populated with several options about how we want to read the data.



Once you have enabled all the options on the data that you are interested in click the **Apply** button to finish loading the data. For a more detailed explanation of the object inspector read the Properties Section

## References

[1] [http://www.paraview.org/Wiki/ParaView/Plugin\\_HowTo#Adding\\_a\\_Reader](http://www.paraview.org/Wiki/ParaView/Plugin_HowTo#Adding_a_Reader)

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# Understanding Data

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## VTK Data Model

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### Introduction

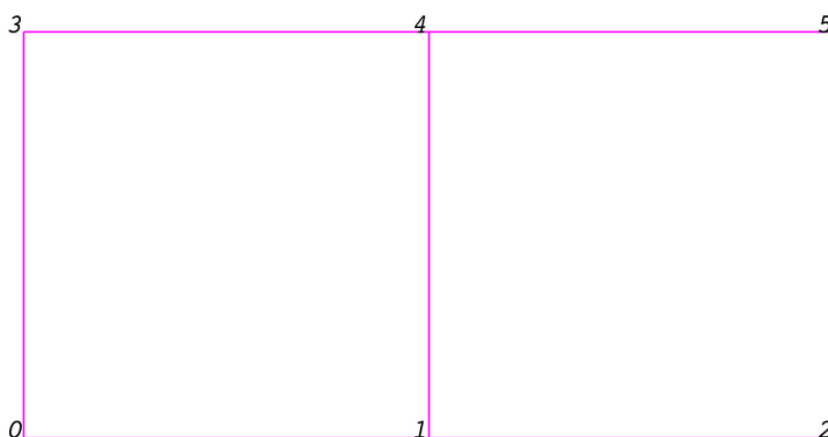
To use ParaView effectively, you need to understand the ParaView data model. Here, we will introduce the VTK data model used by ParaView briefly. For more details, you can refer to one of the VTK books.

The most fundamental data structure in VTK is a data object. Data objects can either be scientific datasets such as rectilinear grids or finite elements meshes (see below) or more abstract data structures such as graphs or trees. Since ParaView is geared toward scientific visualization, we will focus on the former and touch briefly on the latter when necessary.

Datasets are formed of smaller building blocks: mesh (topology and geometry) and attributes.

### Mesh

Even though the actual data structure used to store the mesh in memory depends on the type of the dataset, some abstractions are common to all types. We describe these here and we briefly describe the data structures below for each dataset. In general, a mesh consists of vertices (points) and cells (elements, zones). Cells are used to discretize a region and can have various types such as tetrahedra, hexahedra etc. Each cell contains a set of vertices. The mapping from cells to vertices is called the connectivity. Note that even though it is possible to define data elements such as faces and edges, VTK does not represent these explicitly. Rather, they are implied by a cell's type and its connectivity. One exception to this rule is the arbitrary polyhedron which explicitly stores its faces. Here is an example mesh that consists of 2 cells. The first cell is defined by vertices (0, 1, 3, 4) and the second cell is defined by vertices (1, 2, 4, 5). These cells are neighbors because they share the edge defined by the points (1, 4).

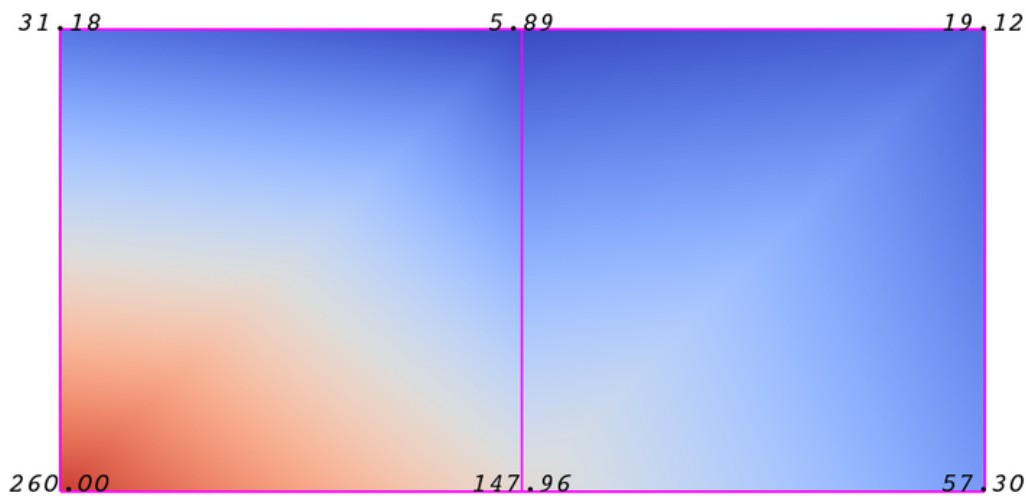


So far, we talked about the topology only. A mesh is fully defined by its topology and the spatial coordinates of its vertices. In VTK, the point coordinates may be implicit or explicitly defined by a data array of dimensions (number\_of\_points, 3).

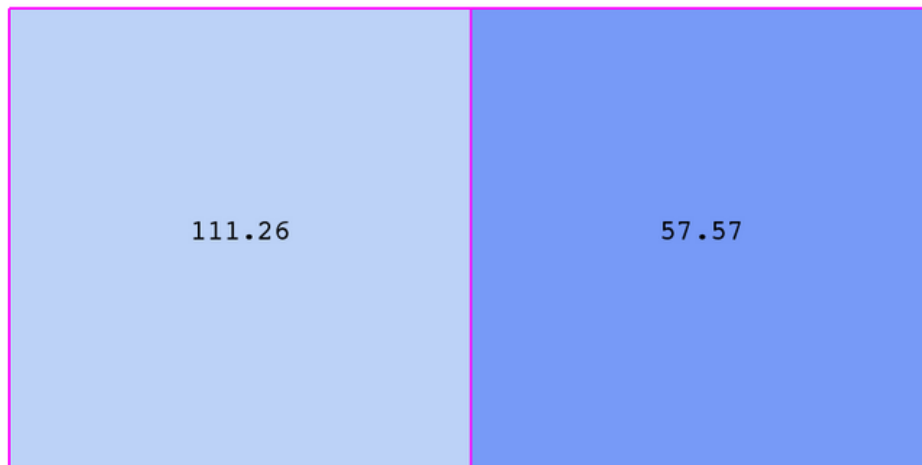
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## Attributes (fields, arrays)

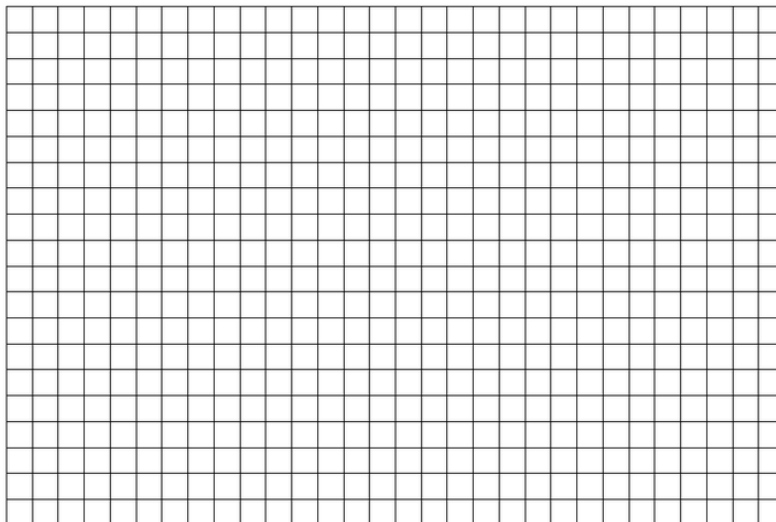
An attribute (aka data array or field) defines the discrete values of a field over the mesh. Examples of attributes include Pressure, Temperature, Velocity and Stress Tensor. Note that VTK does not specifically define different types of attributes. All attributes are stored as data arrays which can have an arbitrary number of components. ParaView makes some assumptions in regards to the number of components. For example, a 3 component array is assumed to be an array of vectors. Attributes can be associated with points or cells. It is also possible to have attributes that are not associated with either. The following example demonstrates the use of a point-centered attribute. Note that the attribute is only defined on the vertices. Interpolation is used to obtain the values everywhere else. The interpolation functions used depends on the cell type. See VTK documentation for details.



The figure below demonstrates the use of a cell-centered attribute. Note that cell-centered attributes are assumed to be constant over each cell. Due to this property, many filters in VTK cannot be directly applied to cell-centered attributes. It is normally required to apply a Cell Data to Point Data filter. In ParaView, this filter is applied automatically when necessary.



## Uniform Rectilinear Grid (Image Data)



A uniform rectilinear grid (aka image data) defines its topology and point coordinates implicitly. To fully define the mesh for an image data, VTK uses the following:

- Extents - these define the minimum and maximum indices in each direction. For example, an image data of extents (0, 9), (0, 19), (0, 29) has 10 points in the x-direction, 20 points in the y-direction and 30 points in the z-direction. The total number of points is  $10 \times 20 \times 30$ .
- Origin - this is the position of a point defined with indices (0, 0, 0)



- Spacing - this is the distance between each point. Spacing for each direction can be defined independently

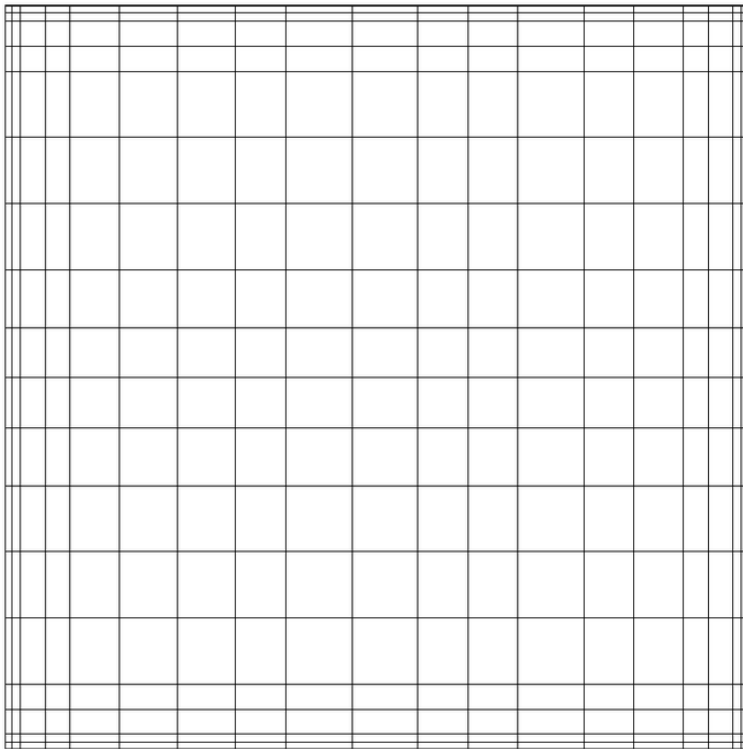
The coordinate of each point is defined as follows:  $coordinate = origin + index * spacing$  where coordinate, origin, index and spacing are vectors of length 3.

Note that the generic VTK interface for all datasets uses a flat index. The (i,j,k) index can be converted to this flat index as follows:  $idx\_flat = k * (npts\_x * npts\_y) + j * npts\_x + i$ .

A uniform rectilinear grid consists of cells of the same type. This type is determined by the dimensionality of the dataset (based on the extents) and can either be vertex (0D), line (1D), pixel (2D) or voxel (3D).

Due to its regular nature, an image data requires less storage than other datasets. Furthermore, many algorithms in VTK have been optimized to take advantage of this property and are more efficient for image data.

## Rectilinear Grid



A rectilinear grid defines its topology implicitly and point coordinates semi-implicitly. To fully define the mesh for a rectilinear grid, VTK uses the following:

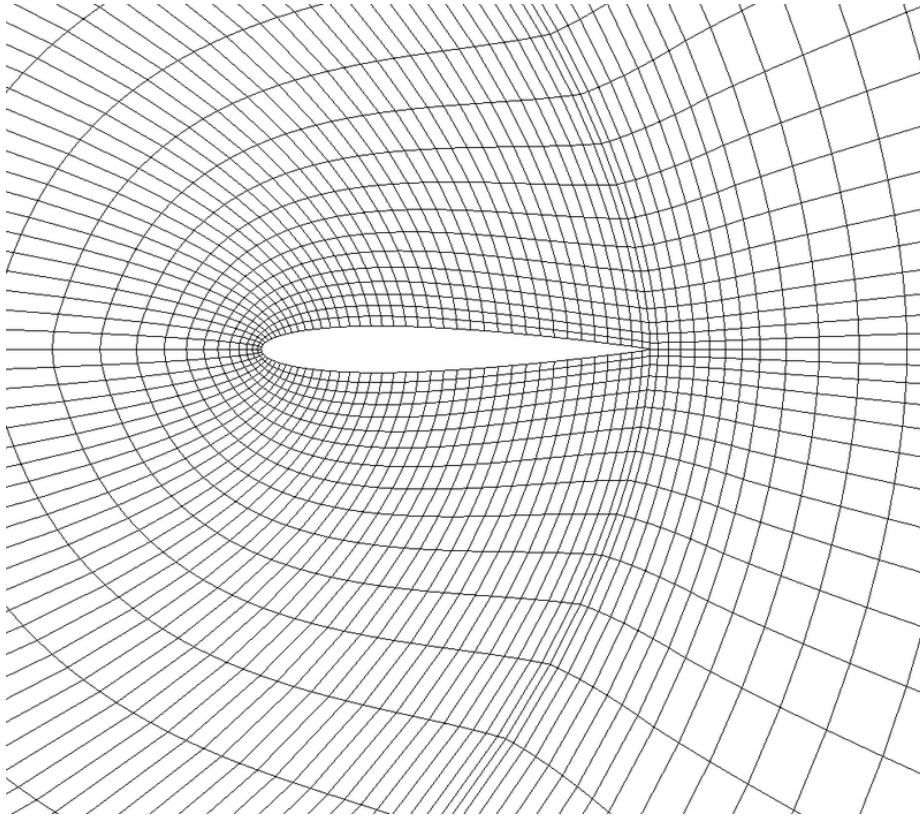
- Extents - these define the minimum and maximum indices in each direction. For example, a rectilinear grid of extents (0, 9), (0, 19), (0, 29) has 10 points in the x-direction, 20 points in the y-direction and 30 points in the z-direction. The total number of points is  $10 * 20 * 30$ .
- 3 arrays defining coordinates in the x-, y- and z-directions. These arrays are of length `npts_x`, `npts_y` and `npts_z`. This is a significant savings in memory as total memory used by these arrays is `npts_x + npts_y + npts_z` rather than `npts_x * npts_y * npts_z`.

The coordinate of each point is defined as follows:  $coordinate = (coordinate\_array\_x(i), coordinate\_array\_y(j), coordinate\_array\_z(k))$ .

Note that the generic VTK interface for all datasets uses a flat index. The (i,j,k) index can be converted to this flat index as follows:  $idx\_flat = k * (npts\_x * npts\_y) + j * npts\_x + i$ .

A rectilinear grid consists of cells of the same type. This type is determined by the dimensionality of the dataset (based on the extents) and can either be vertex (0D), line (1D), pixel (2D) or voxel (3D).

## Curvilinear Grid (Structured Grid)



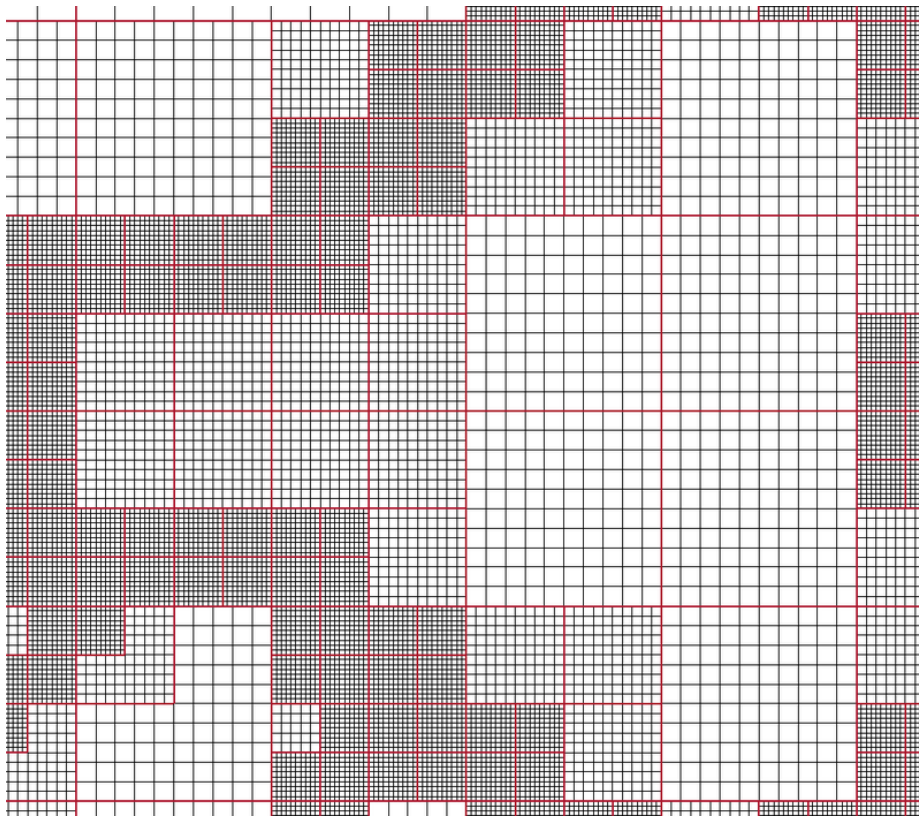
A curvilinear grid defines its topology implicitly and point coordinates explicitly. To fully define the mesh for a curvilinear grid, VTK uses the following:

- Extents - these define the minimum and maximum indices in each direction. For example, a curvilinear grid of extents (0, 9), (0, 19), (0, 29) has 10\*20\*30 points regularly defined over a curvilinear mesh.
- An array of point coordinates. This array stores the position of each vertex explicitly.

The coordinate of each point is defined as follows:  $coordinate = coordinate\_array(idx\_flat)$ . The  $(i, j, k)$  index can be converted to this flat index as follows:  $idx\_flat = k*(npts\_x*npts\_y) + j*nptr\_x + i$ .

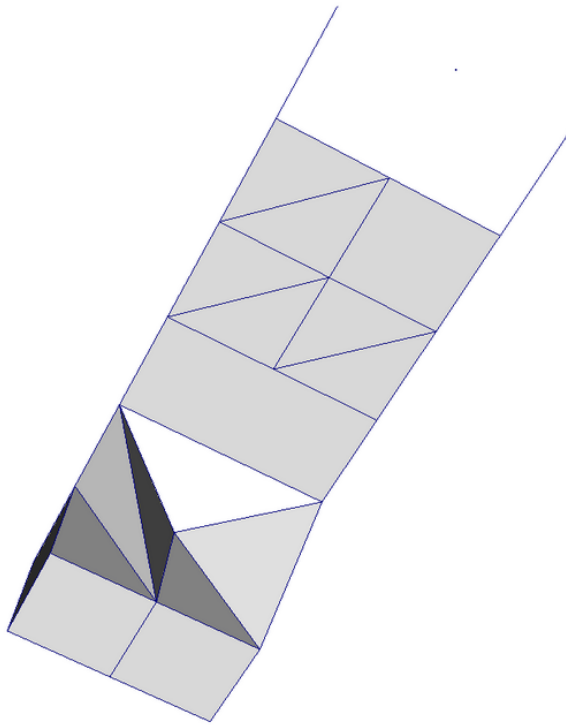
A curvilinear grid consists of cells of the same type. This type is determined by the dimensionality of the dataset (based on the extents) and can either be vertex (0D), line (1D), quad (2D) or hexahedron (3D).

## AMR Dataset



VTK natively support Berger-Oliger type AMR (Adaptive Mesh Refinement) datasets. An AMR dataset is essentially a collection of uniform rectilinear grids grouped under increasing refinement ratios (decreasing spacing). VTK's AMR dataset does not force any constraint on if and how these grids should overlap. However, it provides support for masking (blanking) sub-regions of the rectilinear grids using an array of bytes. This allows VTK to process overlapping grids with minimal artifacts. VTK can automatically generate the masking arrays for Berger-Oliger compliant meshes.

## Unstructured Grid



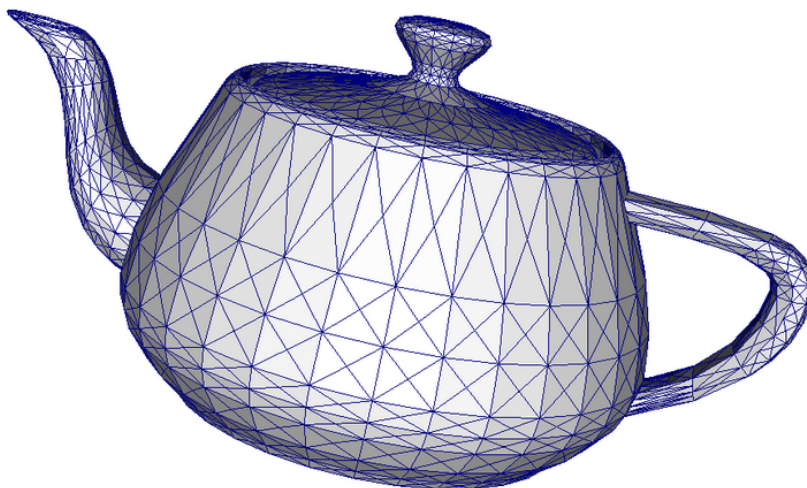
An unstructured grid is most general primitive dataset type. It stores topology and point coordinates explicitly. Even though VTK uses a memory-efficient data structure to store the topology, unstructured grid uses significantly more memory to represent its mesh. Therefore, use an unstructured grid only when you cannot represent your dataset as one of the above datasets. VTK supports a large number of cell types all of which can exist (heterogeneously) within one unstructured grid. The full list of all cell types supported by VTK can be found in the file `vtkCellType.h` in the VTK source code. Here is the list as of when this document was written:

- VTK\_EMPTY\_CELL
- VTK\_VERTEX
- VTK\_POLY\_VERTEX
- VTK\_LINE
- VTK\_POLY\_LINE
- VTK\_TRIANGLE
- VTK\_TRIANGLE\_STRIP
- VTK\_POLYGON
- VTK\_PIXEL
- VTK\_QUAD
- VTK\_TETRA
- VTK\_VOXEL
- VTK\_HEXAHEDRON
- VTK\_WEDGE
- VTK\_PYRAMID
- VTK\_PENTAGONAL\_PRISM
- VTK\_HEXAGONAL\_PRISM
- VTK\_QUADRATIC\_EDGE
- VTK\_QUADRATIC\_TRIANGLE

- VTK\_QUADRATIC\_QUAD
- VTK\_QUADRATIC\_TETRA
- VTK\_QUADRATIC\_HEXAHEDRON
- VTK\_QUADRATIC\_WEDGE
- VTK\_QUADRATIC\_PYRAMID
- VTK\_BIQUADRATIC\_QUAD
- VTK\_TRIQUADRATIC\_HEXAHEDRON
- VTK\_QUADRATIC\_LINEAR\_QUAD
- VTK\_QUADRATIC\_LINEAR\_WEDGE
- VTK\_BIQUADRATIC\_QUADRATIC\_WEDGE
- VTK\_BIQUADRATIC\_QUADRATIC\_HEXAHEDRON
- VTK\_BIQUADRATIC\_TRIANGLE
- VTK\_CUBIC\_LINE
- VTK\_CONVEX\_POINT\_SET
- VTK\_POLYHEDRON
- VTK\_PARAMETRIC\_CURVE
- VTK\_PARAMETRIC\_SURFACE
- VTK\_PARAMETRIC\_TRI\_SURFACE
- VTK\_PARAMETRIC\_QUAD\_SURFACE
- VTK\_PARAMETRIC\_TETRA\_REGION
- VTK\_PARAMETRIC\_HEX\_REGION

Many of these cell types are straightforward. For details, see VTK documentation.

## Polygonal Grid (Polydata)



A polydata is a specialized version of unstructured grid designed for efficient rendering. It consists of 0D cells (vertices and polyvertices), 1D cells (lines and polylines) and 2D cells (polygons and triangle strips). Certain filters

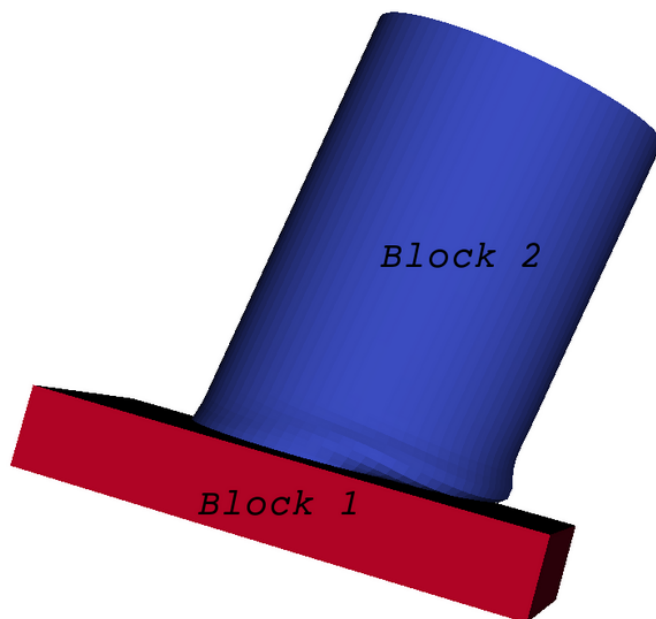
that generate only these cell types will generate a polydata. Examples include the Contour and Slice filters. An unstructured grid, as long as it has only 2D cells supported by polydata, can be converted to a polydata using the Extract Surface filter. A polydata can be converted to an unstructured grid using Clean to Grid.

## Table

|   | Author  | Affiliation |             |     | Alma Mater | Categories  |       | Age | Coolness |
|---|---------|-------------|-------------|-----|------------|-------------|-------|-----|----------|
| 0 | Biff    | NASA        |             |     | Ole...     | Jazz;       | Ro... | 27  | 0.6      |
| 1 | Bob     | Bob's       | Supermarket |     | Ole...     | Jazz        |       | 54  | 0.3      |
| 2 | Baz     | Bob's       | Supermarket |     | TVI        | Food        |       | 16  | 0.3      |
| 3 | Bippity | Oil         | Changes     | 'R' | TVI        | Food        |       | 23  | 0.2      |
| 4 | Boppity | Oil         | Changes     | 'R' | Home       | Food;       | A...  | 34  | 0.25     |
| 5 | Boo     | Oil         | Changes     | 'R' | Princeton  | Automobiles |       | 27  | 0.7      |

A table is a tabular dataset that consists of rows and columns. All chart views have been designed to work with tables. Therefore, all filters that can be shown within the chart views generate tables. Also, tables can be directly loaded using various file formats such as the comma separated values format. Tables can be converted to other datasets as long as they are of the right format. Filters that convert tables include Table to Points and Table to Structured Grid.

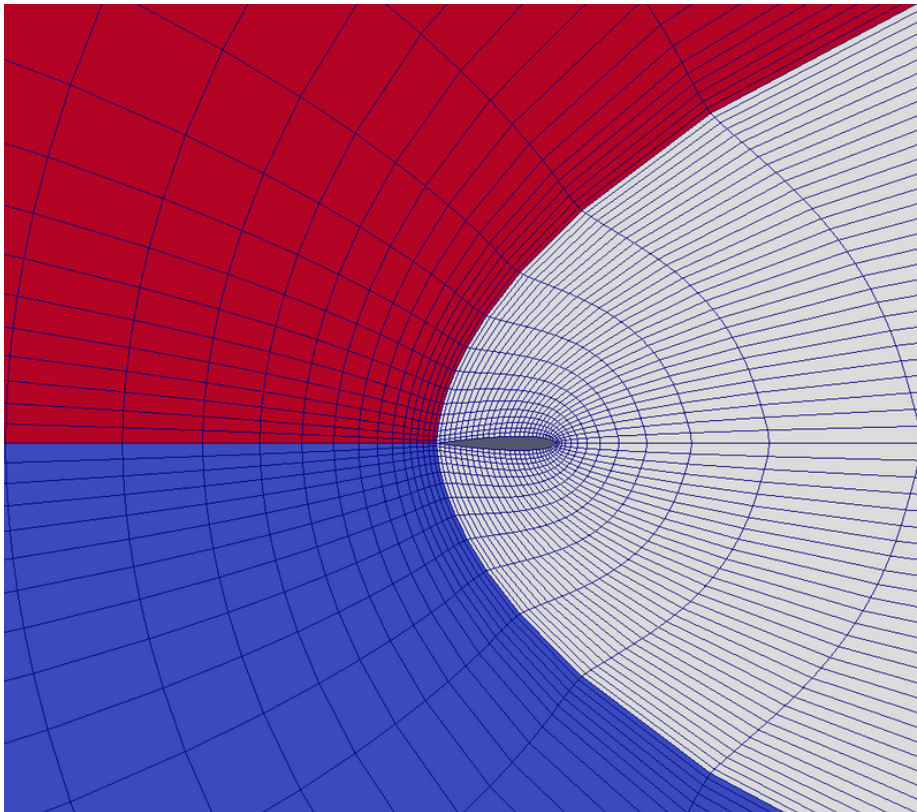
## Multiblock Dataset



You can think of a multi-block dataset as a tree of datasets where the leaf nodes are "simple" datasets. All of the data types describe above, except AMR, are "simple" datasets. Multi-block datasets are used to group together datasets that are related. The relation between these datasets is not necessarily defined by ParaView. A multi-block dataset can represent an assembly of parts or a collection of meshes of different types from a coupled simulation. Multi-block datasets can be loaded or created within ParaView using the Group filter. Note that the leaf nodes of a multi-block dataset don't all have to have the same attributes. If you apply a filter that requires an attribute, it will be

applied only to blocks that have it.

## Multipiece Dataset



Multi-piece datasets are similar to multi-block datasets in that they group together simple datasets. However, there is one big difference. Multi-piece datasets group together datasets that are part of a whole mesh - datasets of the same type and with same attributes. This data structure is used collect datasets produced by a parallel simulation without having to append the meshes together. Note that there is not way to create a multi-piece dataset within ParaView. They are created by certain readers. Furthermore, multi-piece datasets act, for the most part, as simple datasets. For example, it is not possible to extract individual pieces or obtain information about them.

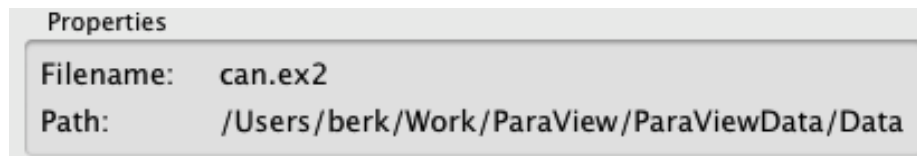


# Information Panel

## Introduction

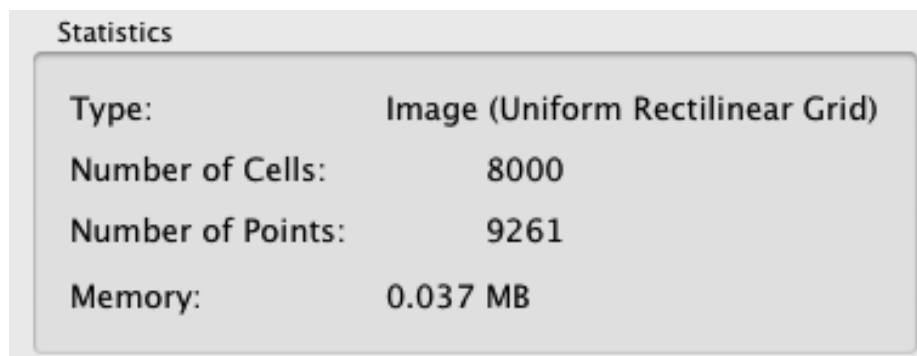
Clicking on the Information button on the Object Inspector will take you to the Information Panel. The purpose of this panel is to provide you with information about the output of the currently selected source, reader or filter. The information on this panel is presented in several sections. We start by describing the sections that are applicable to all dataset types then we describe data specific sections.

## File Properties



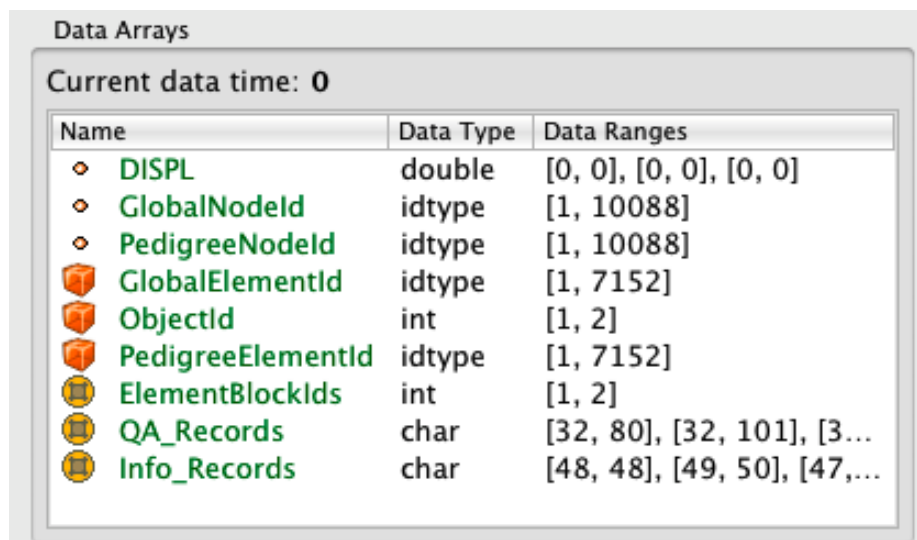
If the current pipeline object is a reader, the top section will display the name of the file and its full path.

## Data Statistics



The Statistics section displays high-level information about the dataset including the type, number of points and cells and the total memory used. Note that the memory is for the dataset only and does not include memory used by the representation (for example, the polygonal mesh that may represent the surface). All of this information is for the current time step.

## Array Information





This section shows the association (point, cell or global), name, type and range of each array in the dataset. In the example, the top 3 attributes are point arrays, the middle 3 cell arrays and the bottom three global (field) arrays. Note that for vectors, the range of each component is shown separately. In case, the range information does not fit the frame, the tooltip will display all of the values.

## Bounds



The Bounds section will display the spatial bounds of the dataset. These are the coordinates of the smallest axis-aligned hexahedron that contains the dataset as well as its dimensions.

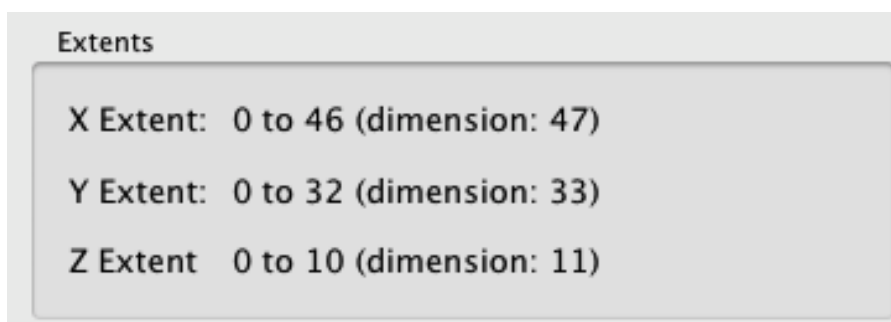
## Timesteps

The screenshot shows a window titled "Time" containing a table with two columns: "Index" and "Value".

| Index | Value       |
|-------|-------------|
| 0     | 0           |
| 1     | 0.000100074 |
| 2     | 0.000199905 |
| 3     | 0.000299964 |
| 4     | 0.000400087 |
| 5     | 0.000499919 |
| 6     | 0.000599935 |
| 7     | 0.000700049 |
| 8     | 0.000800035 |
| 9     | 0.000900061 |

The Time section shows the index and value of all time steps available in a file or produceable by a source. Note that this section display values only when a reader or source is selected even though filters downstream of such sources also have time varying outputs. Also note that usually only one time step is loaded at a time.

## Extents



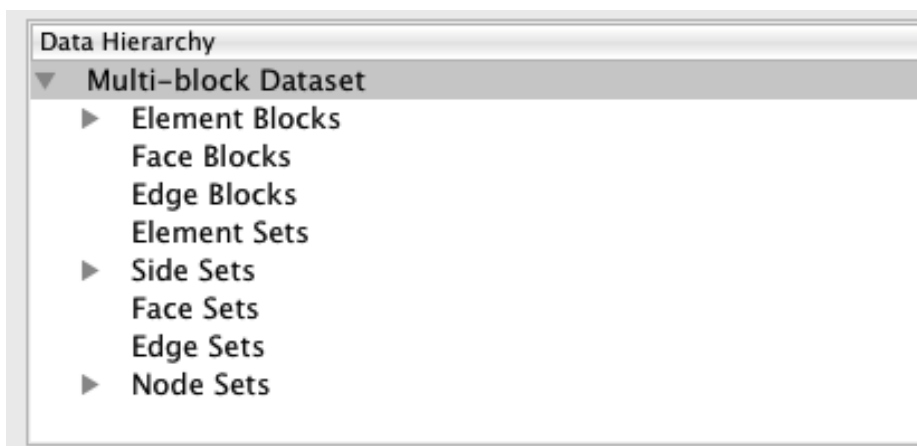
The Extents section is available only for structured datasets (uniform rectilinear grid, rectilinear grid and curvilinear grid). It displays the extent of all three indices that define a structured datasets. It also displays the dimensions (the number of points) in each direction. Note that these refer to logical extents and the labels X Extent, Y Extent and Z Extent can be somehow misleading for curvilinear grids.

## Data Hierarchy (AMR)















For AMR datasets, the Data Hierarchy section shows the various refinement levels available in the dataset. Note that you can drill down to each level by clicking on it. All of the other sections will immediately update for the selected level. For information on the whole dataset, select the top parent called "AMR Dataset".

## Data Hierarchy (Multi-Block Dataset)



For multi-block datasets, the Data Hierarchy section shows the tree that forms the multi-block dataset. By default, only the first level children are shown. You can drill down further by clicking on the small triangle to the left of each node. Note that you can drill down to each block by clicking on it. All of the other sections will immediately update for the selected block. For information on the whole dataset, select the top parent called "Multi-Block Dataset".

# Statistics Inspector

| Statistics Inspector |  |  |  |             |                    |                   |                 |
|----------------------|--|--|--|-------------|--------------------|-------------------|-----------------|
| Name                 | Data Type  | No. of Cells   | No. of Points  | Memory (MB) | Geometry Size (MB) | Spatial Bounds    | Temporal Bounds |
| Wavelet1             |  Image (Uniform Rectilinear Grid) |  8000   |  9261   | 0.037       | 0.002              | [ -10, 10 ] ,...  | [ALL]           |
| RectGrid2.vtk        |  Rectilinear Grid                 |  14720  |  17061  | 0.27        | 0.002              | [ -1.22, 1.1...   | [ALL]           |
| can.ex2              |  Multi-block Dataset              |  7152   |  10088  | 1.626       | 0.795              | [ -7.88, 8.3...   | [ 0, 0.0043]    |
| spcth.0              |  AMR Dataset                      |  103512 |  138999 | 0.114       | 0.071              | [ -3, 5 ] , [ ... | [ 0, 5.01e-07]  |

The Statistics Inspector (View -> Statistics Inspector) can be used to obtain high-level information about the data produced by all sources, readers and filters in the ParaView pipeline. Some of this information (and more) is also available through Object Inspector's Information panel. The information presented in the Statistics Inspector include the name of the pipeline object that produced the data, the data type, the number of cells and points, memory used by the dataset, memory used by the visual representation of the dataset (usually polygonal data), the spatial bounds of the dataset and the temporal bounds of the dataset (the minimum and maximum time values for all available time steps).

Note that the selection in the Statistics Inspector is linked with the Pipeline Browser. Selecting an entry in the Selection Inspector will update the Pipeline Browser and vice versa.

# Displaying Data

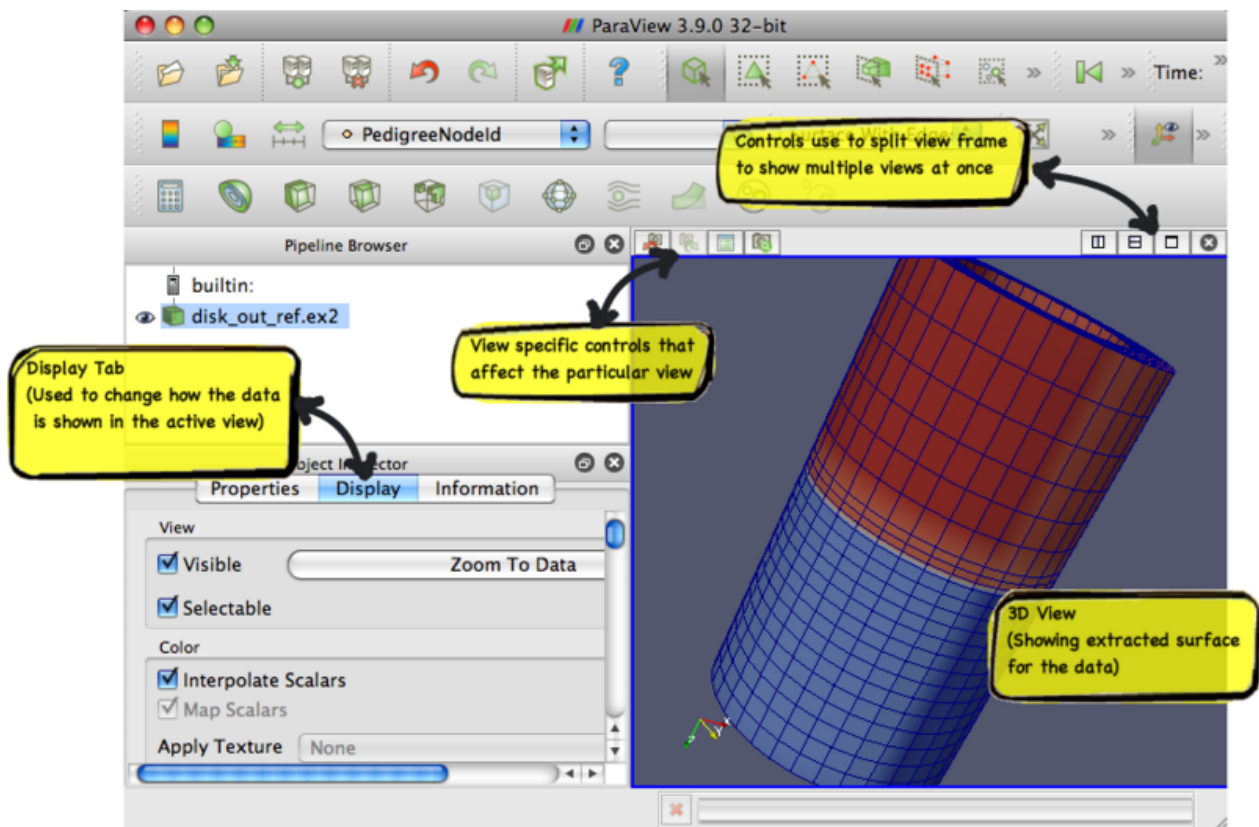
## Views, Representations and Color Mapping

The goal of visualization is to generate representations of the data in a visual form. In this chapter we will see different mechanisms available in ParaView to look at the data to gain insight into it.

### Understanding Views

#### Views

When the ParaView application starts up, one sees a 3D viewport with an axes at the center. This is a *View*. In ParaView, *Views* are frames in which the data can be *seen*. There are different types of views. The default view that shows up is a *3D View* which shows rendering of the geometry extracted from the data or volumes or slices in a 3D scene. You can change the default view **Settings** dialog (Edit | Settings (in case of Mac OS X, ParaView | Preferences)).

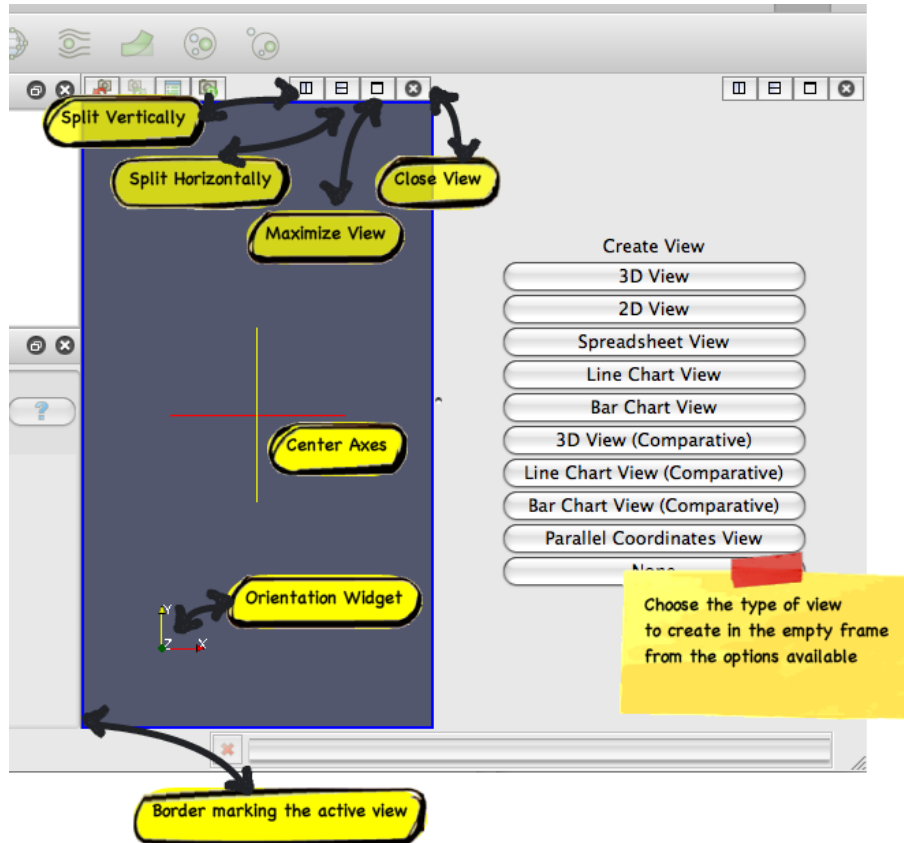


There may be parameters that are available to the user that control how the data is displayed e.g. in case of 3D View, the data can be displayed as wireframes or surfaces, you choose to change the color of the surface or use a scalar for coloring etc. All these options are known as *Display* properties and are accessible from the **Display** tab in the **Object Inspector**.

Since there can be multiple datasets shown in a view, as well as multiple views, the **Display** tabs show the properties for the active pipeline object (changed by using the Pipeline browser, for example) in the active view.

## Multiple Views

ParaView supports showing multiple views side by side. To create multiple views, simply use the controls on the top-right corner of the view to split the frame vertically or horizontally. You can even maximize a particular view to temporarily hide other views. Once a view-frame is split, you will see a list of buttons showing the different types of views that you can create to place in that view. Simply click the button to create the view of your choice.



Some filters, such as **Plot Over Line** may automatically split the view frame and show the data in a particular type of view suitable for the data generated by the filter.

## Active View

Once you have multiple views, the active view is indicated by a colored border around the view frame. Several menus as well as toolbar buttons affect the active view alone. Also they may become enabled/disabled based on whether that corresponding action is supported by the active view.

The **Display** tab affects the active view. Similarly, the *eye icon* in the **Pipeline Browser**, next to the pipeline objects, indicates the visibility state for that object in the active view.

When a new filter or source or reader is created, it will be displayed by default in the active view, if possible (otherwise, it may create a new view).

## Types of Views

In this section we will cover the different types of Views available in ParaView. For each view, we will talk about the controls available to change the view parameters using **View Settings** as well as the parameters associated with the **Display Tab** for showing data in that view.

### 3D View

3D View is used to show the surface or volume rendering for the data in a 3D world. This is the most commonly used view type.

When running in client-server mode, 3D View can render data either by bringing the geometry to the client and then rendering it there or by rendering it on the server (possibly in parallel) and then delivering the composited images to the client. Refer to the **Client-Server Visualization** chapter for details.

### Interaction

Interacting with the 3D view will typically update the camera. This makes it possible to explore the visualization scene. The default buttons are as follows, they can be changed using the **Application Settings** dialog.

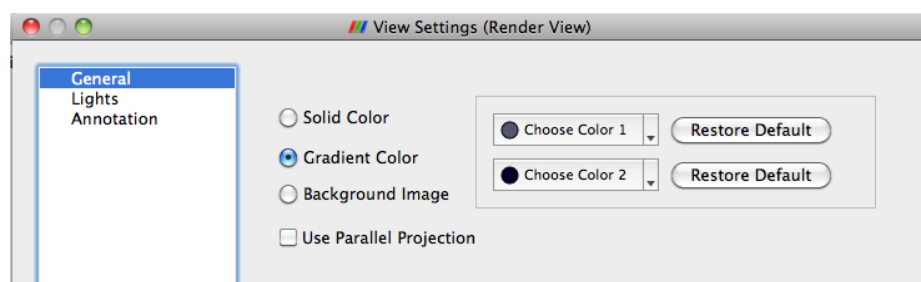
| Modifier | Left Button | Middle Button | Right Button |
|----------|-------------|---------------|--------------|
|          | Rotate      | Pan           | Zoom         |
| Shift    | Roll        | Rotate        | Pan          |
| Control  | Zoom        | Rotate        | Zoom         |

This view supports selection. One can select cells or points either on the surface or those within a frustum. Selecting cells or points makes it possible to extract those for further inspection or label them etc. Details about data querying and selection can be found the Quantitative analysis chapter.

### View Settings

The view settings dialog accessible through the Edit | View Settings menu or the tool button on the left-corner of the view is used to change the view settings per view.

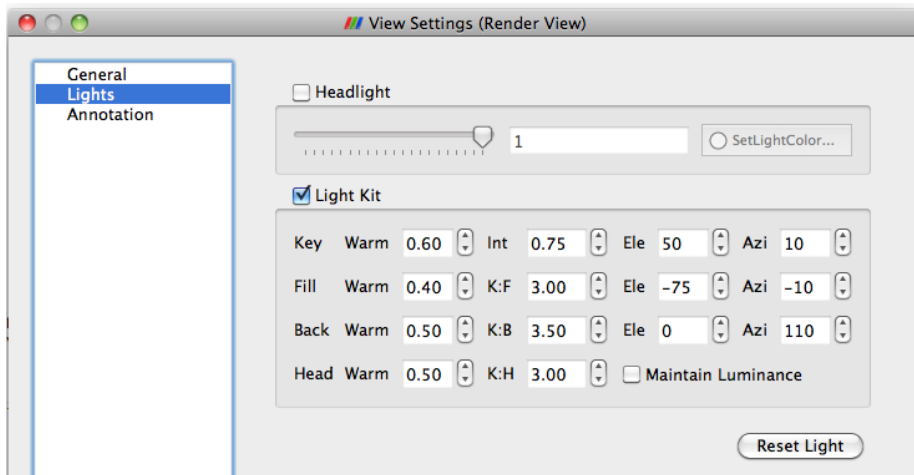
#### General



This tab allows the user to choose the background color. One can use a solid color or a gradient or even a background image.

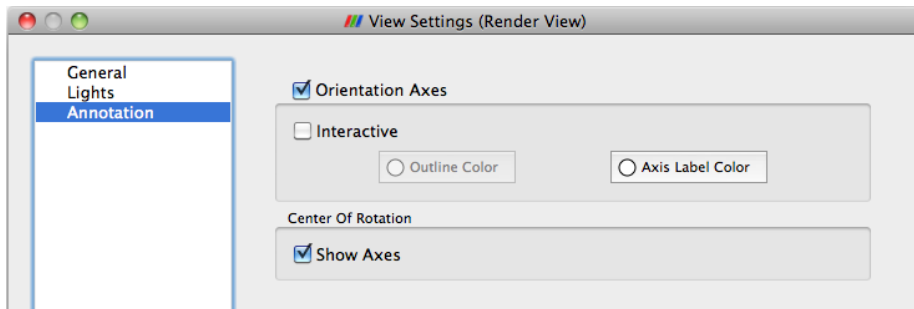
By default the camera uses perspective projection. To switch to parallel projection, check the "Use Parallel Projection" checkbox in this panel.

## Lights



The 3D View requires lights to illuminate the geometry being rendered in the scene. Once can control these lights using this pane.

## Annotation

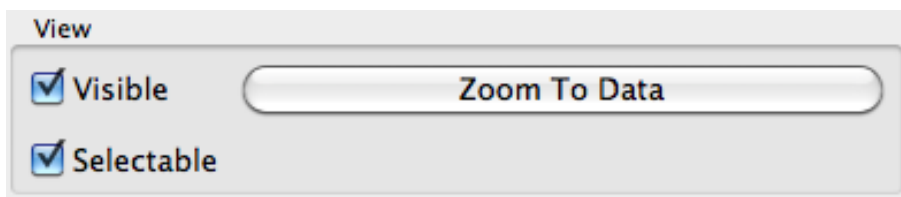


The annotation pane enables controlling the visibility of the center axes and the orientation widget. Users can also make the orientation widget interactive, so that they can manually place the widget at location of their liking.

## Display Properties

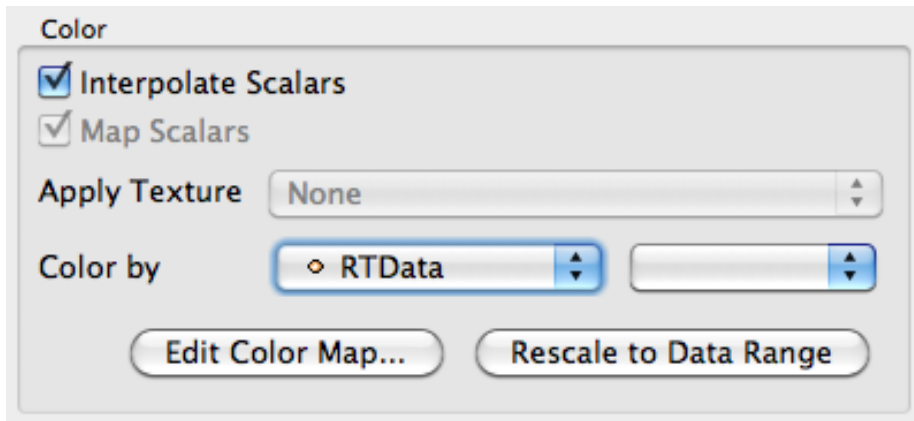
Users can control how the data from any source or filter is shown in this view using the **Display** tab. In this section we will cover the various options available to a user for controlling appearance of the rendering in the 3D view.

## View



| Name         | Usage  |
|--------------|--|
| Visible      | Checkbox used to toggle the visibility of the data in the view. If it disabled, it implies that the data cannot be shown in this view. |
| Selectable   | Checkbox used to toggle whether the data gets selected when using the selection mechanism for selecting and sub-setting data.          |
| Zoom to Data | Click this button to zoom the camera so that the dataset is completely fits within the viewport.                                       |

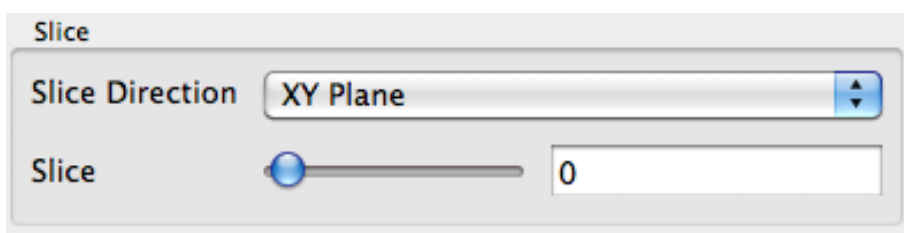
## Color



The color group allows users to pick the scalar to color with or set a fixed solid color for the rendering.

| Name                | Usage  |
|---------------------|--|
| Interpolate Scalars | If on, the scalars will be interpolated within polygons and the scalar mapping happens on per pixel basis. If off, then color mapping happens at points and colors are interpolated which is typically less accurate. This only affects when coloring with point arrays. Has no effect otherwise. This is disabled when coloring using a solid color.  |
| Map Scalars         | If the data array being color with can be directly interpreted as colors then one can uncheck this to not use any lookup table. Otherwise, when on, a lookup table will be used to map scalars to colors. This is disabled when the array is not of the type that can be interpreted as colors (i.e. vtkUnsignedCharArray).  |
| Apply Texture       | Makes is possible to apply a texture over the surface. This requires that the data has texture coordinates. One can use filters like <b>Texture Map to Sphere</b> , <b>Texture Map to Cylinder</b> or <b>Texture Map to Plane</b> to generate texture coordinates when not present in the data. To load a texture, select <b>Load</b> from the combo-box which will popup a dialog allowing you to choose an image. Otherwise pick from already loaded textures listed in the combo-box.   |
| Color By            | This enables coloring of the surface/volume. Either choose the array to color with or set the solid color to use. When volume rendering, solid coloring is not possible, one has to choose the data array to volume render with.   |
| Set solid color     | Used to set the solid color. This is available only when <b>Color By</b> is set to use <i>Solid Color</i> . ParaView defines a notion of a color palette consisting of different color categories. To choose a color from one of these predefined categories, click on the tiny arrow next to this button. It will open up a drop down with options to choose from. If one uses a color from the palette, it possibly to globally change the color by changing the color palette e.g. for printing or for display on screen etc. |
| Edit Color Map...   | One can edit the color-map or lookup table by clicking the Edit Color Map button. It's only shown when an array is chosen in the <b>Color By</b> combo-box.  |

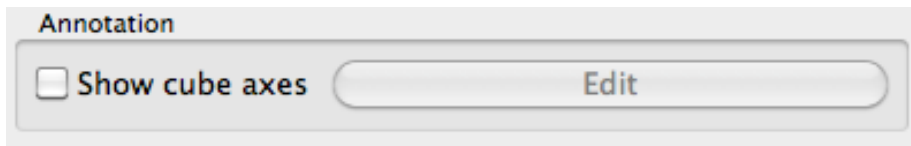
## Slice



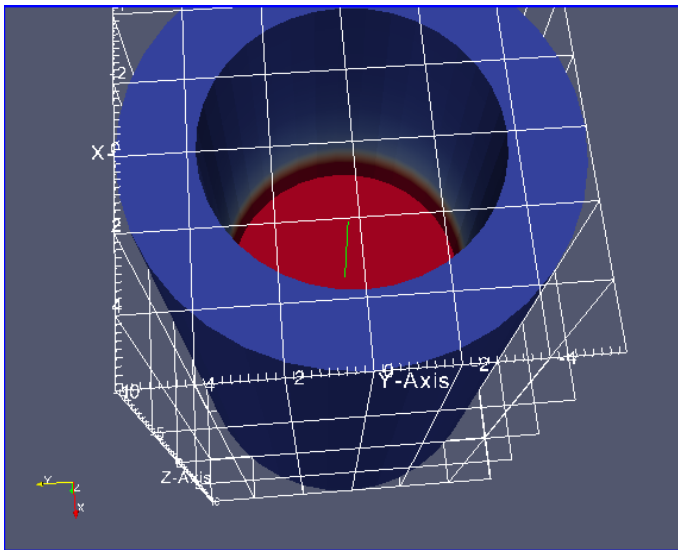


The slice controls are available only for Image dataset (Uniform Rectilinear Grids) when the representation type is **Slice**. The representation type is controlled using the *Style* group on the **Display** tab. These allow the user to pick the slice direction as well as the slice offset.

### Annotation



Cube axes is a annotation box that can be used so show a scale around the dataset. Use the **Show cube axes** checkbox to toggle its visibility. You can further control the apperance of the cube axes by clicking **Edit** once the cube-axes is visible.



Style

Style

RepresentationSurface

InterpolationGouraud

MaterialNone

Point size2.00

Line width1.00

Opacity1.00

Subdivision0

Volume mapperFixed point

Edge Style

Set Edge Color...

| Name           | Usage  |
|----------------|--|
| Representation | Use this to change how the data is represented i.e. as a surface, volume, wireframe, points, or surface with edges.  |
| Interpolation  | Choose the method used to shade the geometry and interpolate point attributes.   |
| Point Size     | If your dataset contains points/vertices, this adjust the diameter of the rendered points. It also affects the point size when <b>Representation</b> is <i>Points</i> .  |
| Line width     | If your dataset contains lines/edges, this scale adjust the width of the rendered lines. It also affects the rendered line width when <b>Representation</b> is <i>Wireframe</i> or <i>Surface With Edges</i> . |
| Opacity        | Set the opacity of the dataset's geometry. ParaView uses hardware-assisted Depth peeling, whenever possible, to remove artifacts due incorrect sorting order of rendered primitives.                           |
| Volume Mapper  | When <b>Representation</b> is <i>Volume</i> , this combo-box allows the user to choose a specific volume rendering technique. The techniques available change based on the type of the dataset.                |
| Set Edge Color | This is available when <b>Representation</b> is <i>Surface with Edgets</i> . It allows the user to pick the color to use for the edges rendered over the surface.  |

Backface Style

Backface Style

RepresentationCull Frontface

Set Backface Color...

Opacity1.00

This allows the user to define backface properties. In computer graphics, backface refers to the face of a geometric primitive with the normal point away from the camera. Users can choose to hide the backface or front face or specify different characteristics for the two faces using these settings.

### Transformation

| Transformation |   |   |   |
|----------------|---|---|---|
| Translate      | 0 | 0 | 0 |
| Scale          | 1 | 1 | 1 |
| Orientation    | 0 | 0 | 0 |
| Origin         | 0 | 0 | 0 |

These settings allow the user to transform the rendered geometry, without actually transforming the data. Note that since this transformation happens during rendering, any filters that you apply to this data-source, will still be working on the original, untransformed data. Use the **Transform** filter if you want to transform the data instead.

### 2D View

2D View is similar to 3D view except that it can only show slices from uniform grid datasets i.e. volume datasets or image datasets. The interaction is also limited to a 2D plane i.e. rotation is not available. Currently this view does not support selection. That will be supported in future releases. When one loads a 2D image, ParaView by default creates this view.

Since this view is same as the 3D view, it has the same features when it comes to client-server and parallel rendering.

### Interaction

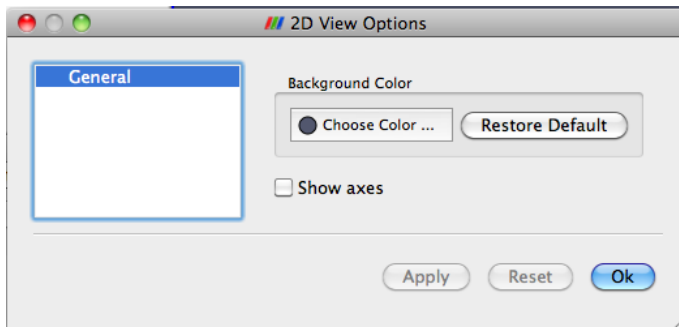
Interactions with this view update the camera. The default buttons are as follows, they can be changed using the **Application Settings** dialog.

| Modifier | Left Button | Middle Button | Right Button |
|----------|-------------|---------------|--------------|
|          | Pan         | Pan           | Zoom         |
| Shift    | Zoom        | Zoom          | Zoom         |
| Control  | Zoom        | Zoom          | Pan          |

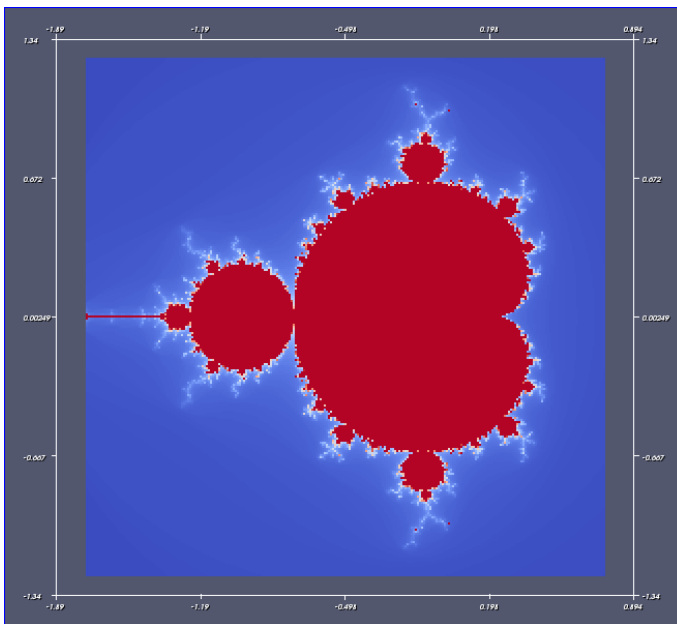
Unlike 3D view, this view currently does not support selection. However, that support will be added in future.

## View Settings

The view settings dialog accessible through the Edit | View Settings menu or the tool button on the left-corner of the view is used to change the view settings per view.



The user can change the background color and toggle the visibility of the axes-box.



## Display Properties

Display tab for this view is exactly same as the 3D View, except that some options may be unavailable such as changing of representation type.

## Spreadsheet View

Spreadsheet View is used to inspect the raw data in a spreadsheet. When running in client-server mode, to avoid delivering the entire dataset to the client for displaying the spreadsheet (since the data can be very large), this view streams only visible chunks of the data to the client. As the user scrolls around the spreadsheet, new data chunks are fetched.

Unlike some other views, this view can only show one dataset at a time. For composite datasets, it shows only one block at a time. One can pick the block to show using the **Display** tab.

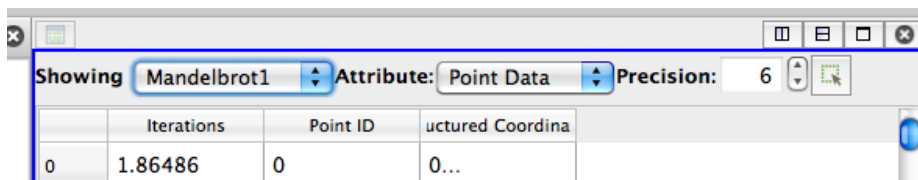
## Interaction

As far as the usability is concerned, this view behaves exactly like typical spreadsheets shown in applications like Microsoft's Excel or Apple's Pages:

- One can scroll up and down to inspect new rows.
- One can sort any column by clicking on the header for the column. Repeated clicking on the column header toggles the sorting order. When running in parallel, ParaView uses sophisticated parallel sorting algorithms to avoid memory and communication overheads to sort large, distributed datasets.
- One can double-click on a column header to toggle a mode in which only that column is visible. This reduces clutter when one is interested in a single attribute array.
- One can click on rows to select the corresponding elements i.e. cells or points. This is not available when in "Show selected only mode" (explained later). Also, when one creates a selection in other views e.g. the 3D view, the rows corresponding to the selected elements will be highlighted.

## Header

Unlike other views, Spreadsheet View has a header. This header provides quick access to some of the commonly used functionality in this view.

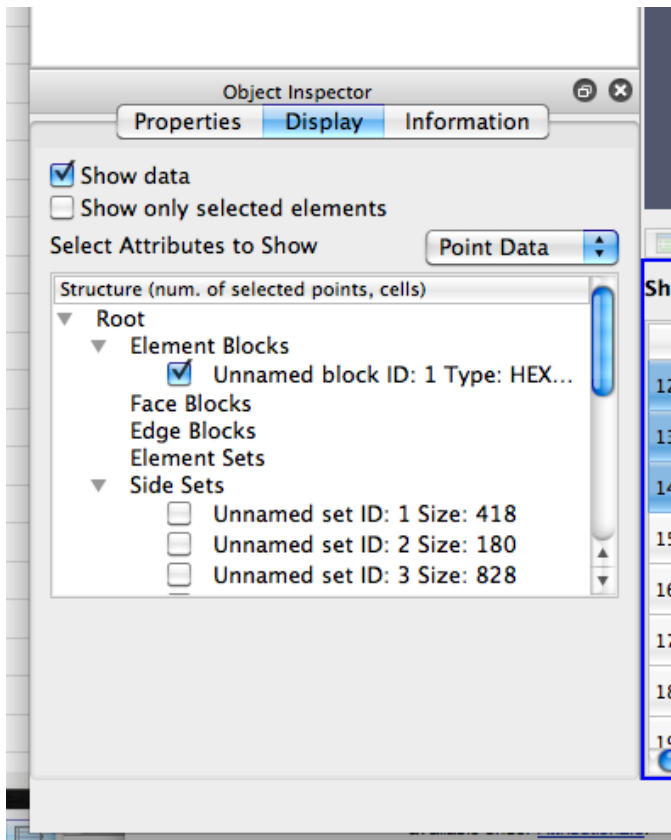


- Since this view can only show one dataset at a time, one can quickly choose the dataset to show using the **Showing** combo-box.
- One can choose the attribute type i.e. point attributes, cell attributes, to display using the **Attribute** combo-box.
- **Precision** controls the number of digits to show after decimal point for floating point numbers.
- The last push-button allows the user to enter the view in a mode where it only shows the selected rows. This is useful when one creates a selection using another view such as the 3D view and wants to inspect the details for the selected cells/points.

## View Settings

Currently, no user settable settings are available for this view.

## Display Properties



The display properties for this view provide the same functionality as the header. Additionally, when dealing with composite datasets, the display tab shows a widget allowing the user to choose the block to display in the view.

## Line Chart View

A traditional 2D line plot is often the best option to show trends in small quantities of data. A line plot is also a good choice to examine relationships between different data values that vary over the same domain.

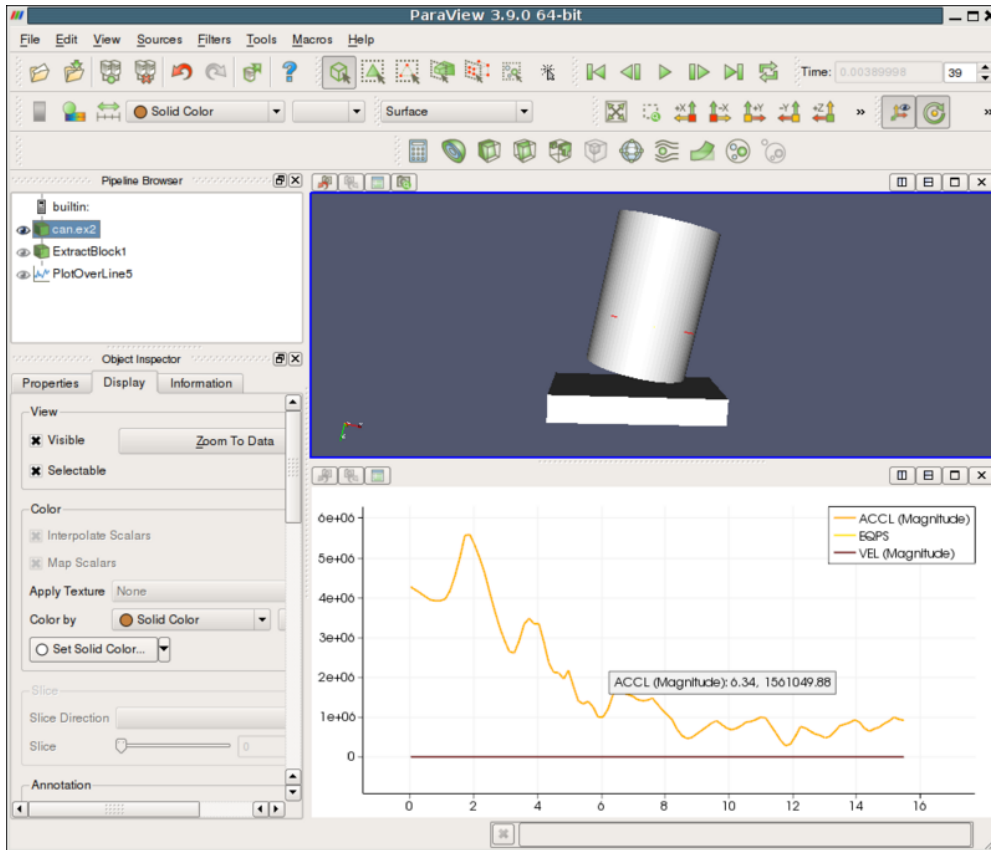
Any reader, source, or filter that produces plottable data can be displayed in an XY Plot View. ParaView stores its plottable data in table (vtkTable). Using the display properties, users can choose which columns in the table must be plotted on the x and y axes.

As with the other view types, exactly what is displayed in the active XY Plot View is displayed by, and controllable with, the eye icons in the Pipeline Browser panel. When an XY Plot View is active, only those filters that produce plottable output have eye icons.

The XY Plot View is the preferred View type for the Plot over Line, Plot Point over Time, Plot Cell over Time, Plot Field Variable over Time, and Probe Location over Time filters. Creating any one of these filters will automatically create an XY Plot View for displaying its output. Figure 3 shows a plot of the data values within a volume as they vary along three separate paths. The top curve comes from the line running across the center of the volume, where the largest values lie. The other two curves come from lines running near the edges of the volume.

Unlike the 3D and 2D render view, the charting views are client-side views i.e. they deliver the data to be plotted to the client. Hence ParaView only allows results from some standard filters such as **Plot over Line** in the line chart view by default. However it is also possible to plot cell or point data arrays for any dataset by apply the **Plot Data**

filter.



## Interaction

Line chart view supports the following interaction modes:

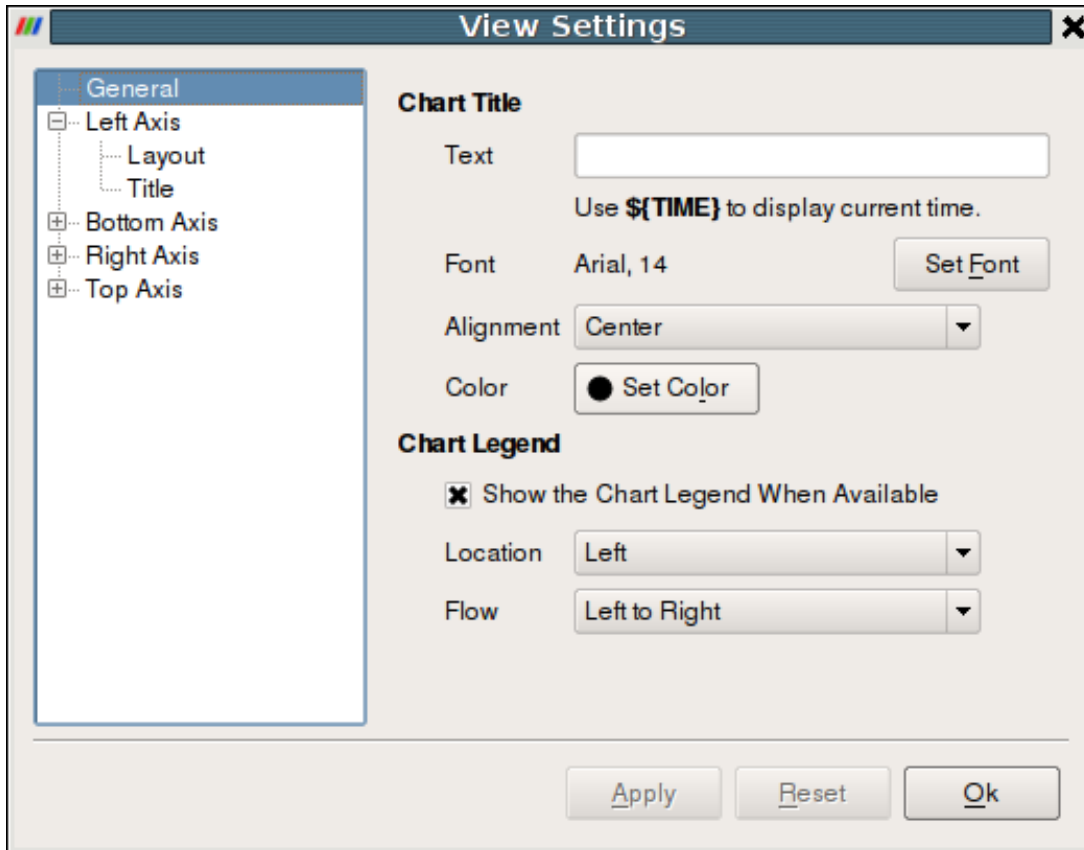
- Right-click and drag to pan
- Left-click and drag to select
- Middle-click and drag to zoom to region drawn.
- Hover over any line in the plot to see the details for the data at that location.

To reset the view, use the **Reset Camera** button in the **Camera Toolbar**.

## View Settings

The **View Settings** for Line Chart enable the user to control the appearance of the chart including titles, axes positions etc. There are several pages available in this dialog. The **General** page controls the overall appearance of the chart, while the other pages controls the appearance of each of the axes.

### General Settings Page



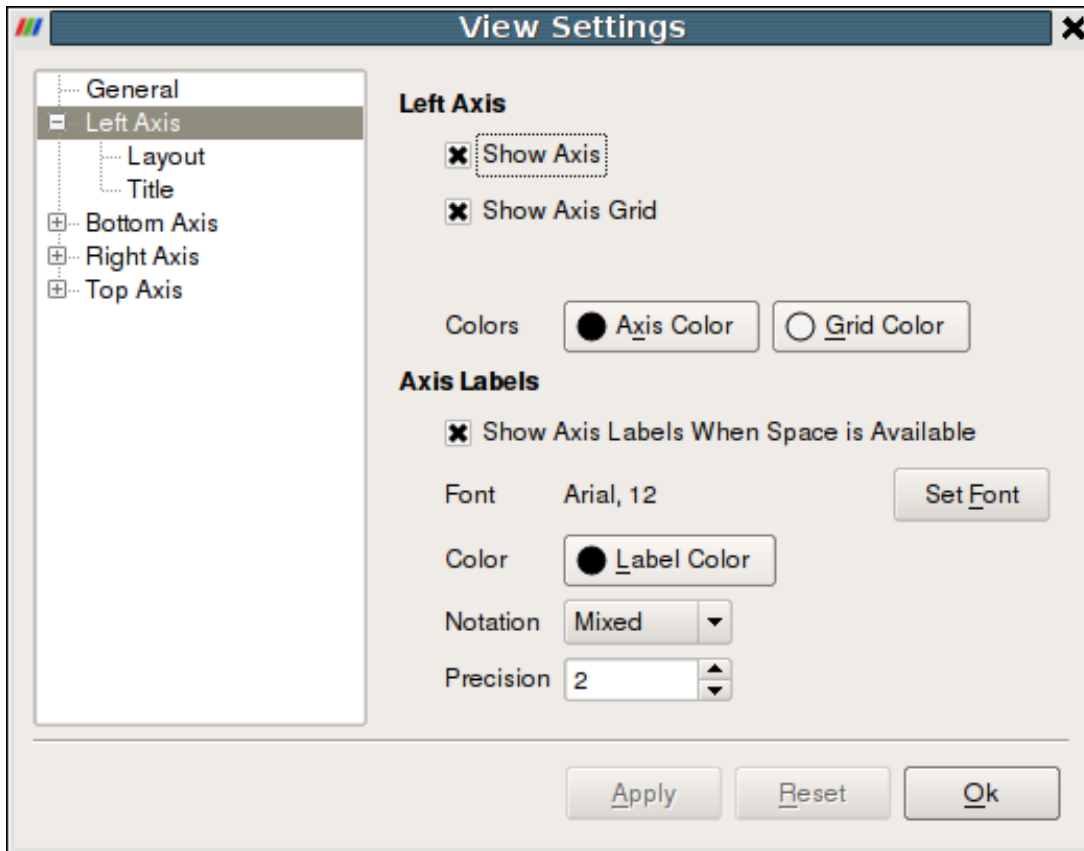
This page allows users to change the title and legend. One can change the title text. To show the current animation time in the title text, simply use the keyword `${TIME}`. Users can further change the font and alignment for the title.

This page also enable changing the appearance and positioning of the legend.



### Axis Settings Page

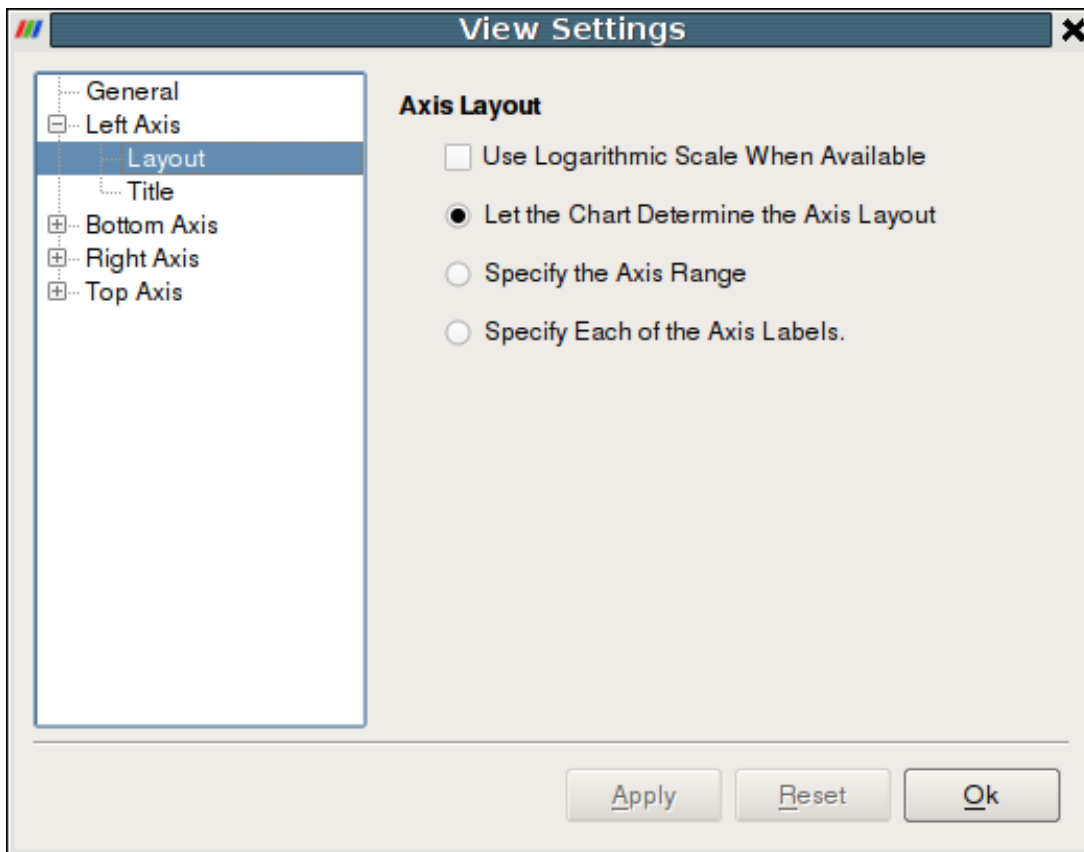
This page allows changing the properties of a particular axis. Four pages are provided for each of the axes. By clicking on the name of the axis, one can access the settings page for the corresponding axes.



- Axis
  - **Show Axis** : controls the axis visibility.
  - **Show Axis Grid**: controls whether a grid is to be drawn perpendicular to this axis.
  - **Colors**: enables changing the axis as well as the grid color.
- Labels
  - **Show Axis Labels When Space is Available** : controls label visibility along this axis.
  - **Font** and **Color**: enables changing the label font and color.
  - **Notation**: allows user to choose between **Mixed**, **Scientific** and **Fixed** point notations for numbers.
  - **Precision**: controls precision after '.' in **Scientific** and **Fixed** notations.

### Axis Layout Page

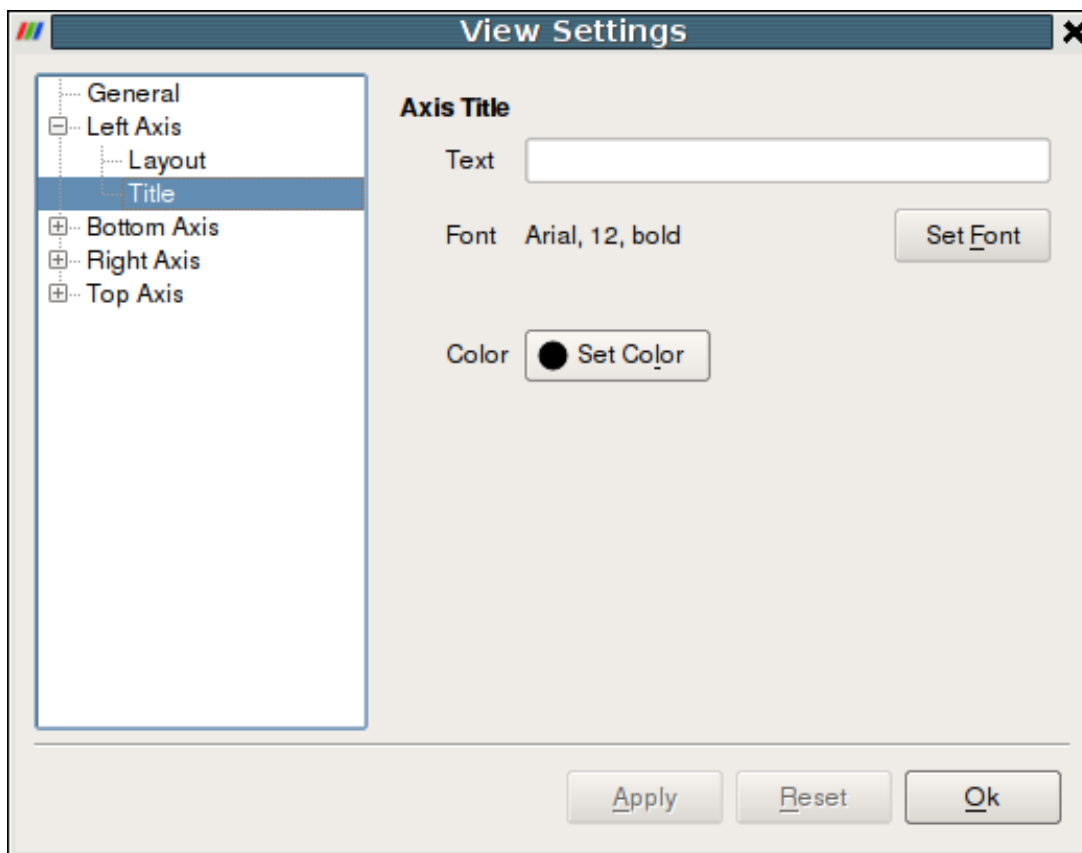
This page allows the user to change the axis range and scale.



- **Use Logarithmic Scale When Available:** check this to use a log scale unless the data contains numbers  $\leq 0$ .
- **Let the Chart Determine the Axis Layout:** pick this radio button to let the chart use the optimal range and spacing for this axis.
- **Specify the Axis Range:** pick this radio button to specify the axis range explicitly. The labels are auto-placed within this range.
- **Specify Each of the Axis Labels:** pick this radio button to specify each label location on axis explicitly.

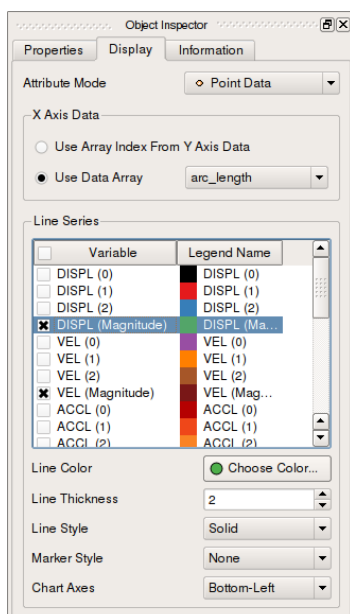
## Axis Title Page

This page allows the user to change the title for the axis. One can change the text as well as color and font for the axis title.



## Display Properties

Display Properties for the Line Chart view allow the user to choose what arrays are plotted along which of the axes as well as the appearance for each of the lines such as its color, thickness and style.



- **Attribute Mode** : pick which attribute arrays to plot i.e. point arrays or cell arrays etc.
- **X Axis Data** : controls the array to use as the X-axis.

- **Use Array Index From Y Axis Data** : when checked results in ParaView simply using the index in data-array plotted on Y as the x axis.
- **Use Data Array**: when checked the user can pick an array to be interpreted as the x co-ordinate.
- **Line Series** : controls the properties of each of the arrays plotted along the y-axis.
  - **Variable** : check the variable to be plotted.
  - **Legend Name** : click to change the name used in the legend for this array.

Select any of the series in the list to change following properties for that series. One can select multiple entries to change multiple series.

- **Line Color** : controls the color for the series.
- **Line Thickness** : controls the thickness for the series.
- **Line Style** : controls the style for the line.
- **Marker Style** : markers can be placed at every data point. This controls the style used for those markers.
- **Chart Axes** :

## Bar Chart View

Traditional 2D graphs present some types of information much more readily than 3D renderings do; they are usually the best choice for displaying one and two dimensional data. The Bar Chart View is very useful for examining the relative quantities of different values within data, for example.

The Bar Chart View is used most frequently to display the output of the Histogram filter. This filter divides the range of a component of a specified array from the input data set into a specified number of bins, producing a simple sequence of the number of values in the range of each bin. A Bar Chart is the natural choice for displaying this type of data. In fact the Bar Chart View is the preferred View type for the Histogram filter. Filters that have a preferred View type will create a View of the preferred type whenever they are instantiated.

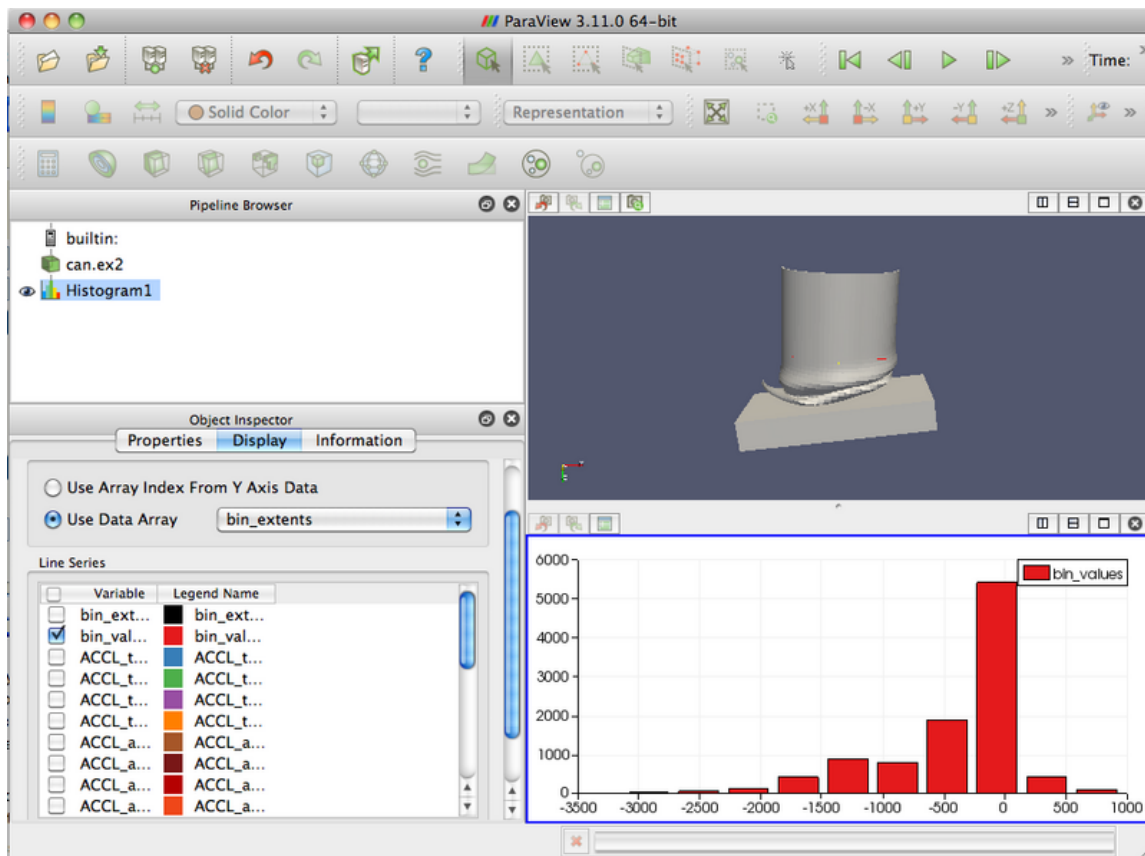
When the new View is created for the histogram filter, the pre-existing 3D view is made smaller to make space for the new Chart View. The Chart View then becomes the Active View, which is denoted with a red border around the View in the display area. Clicking on any View window makes it the active View. The contents of the Object Inspector and Pipeline Browser panels change and menu items are enabled or disabled whenever a different View becomes active to reflect the Active View's settings and available controls. In this way, you can independently control numerous Views. Simply make a View active, and then use the rest of the GUI to change it. By default, the changes you make will only affect the active View.

As with the 3D View, the visibility of different datasets within a Bar Chart View is displayed and controlled by the eye icons in the Pipeline Browser. The Bar Chart View can only display datasets that contain chartable data, and when a Bar Chart View is active, the Pipeline Browser will only display the eye icon next to those datasets that can be charted.

ParaView stores its chartable data in 1D Rectilinear Grids, where the X locations of the grid contain the bin boundaries, and the cell data contain the counts within each bin. Any source or filter that produces data in this format can be displayed in the Bar Chart View. Figure 46 shows a histogram of the values from a slice of a data set.

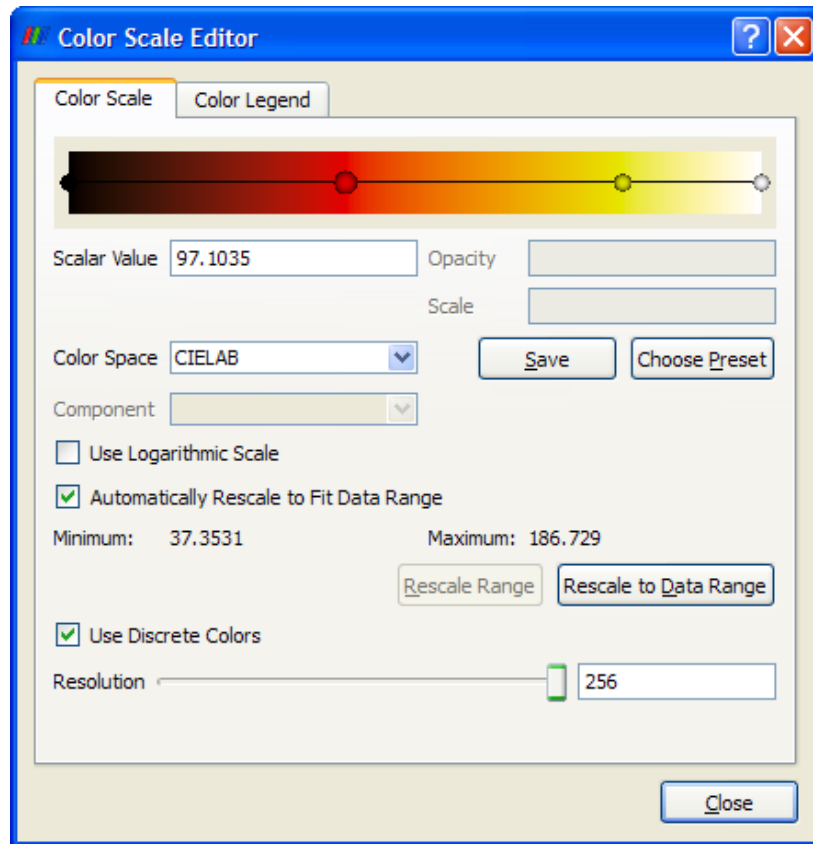
The Chart View's Edit View Options dialog allows you to create labels, titles, and legends for the chart and to control the range and scaling of each axis.

The **Interaction**, **Display Properties** as well as **View Settings** for this view and similar to those for the Line Chart.



## Color Transfer Functions

The interface for changing the color mapping and properties of the scalar bar is accessible from the **Display** tab of the **Object Inspector**. Pressing the **Edit Color Map** button displays the interface for manipulating the color map and scalar bar as shown below.



**Figure 1. Color Map Controls**

At the top of the **Color Scale** tab of the **Color Scale Editor** dialog is an editor for mapping scalar values to colors (for surface rendering) and also to opacity (for volume rendering). The scalar range of this color map editor is shown below the **Automatically Rescale to Fit Data Range** check box. The leftmost sphere corresponds to the minimum scalar value, and the rightmost one corresponds to the maximum. Any interior nodes correspond to values between these two extremes. New nodes may be added to the color editor by left-clicking in the editor; this determines the scalar value associated with the node, but that value may be changed by typing a new value in the **Scalar Value** text box below the color map editor or by clicking and dragging a node. (The scalar value for a particular node may not be changed such that it is less than that for a node left of it or greater than that for a node right of it.) When volume rendering, the vertical height of a node indicates its opacity. For surface rendering, opacity is determined for an entire data set, not based on the underlying scalar values.

When a node in the color map editor is clicked, it becomes highlighted (i.e., drawn larger than the other spheres in the editor). In the example above, the second node from the left has been selected. Clicking again on the selected node displays a color chooser from which you may select a new color for the node. Pressing the 'd' or Delete key while a node is selected removes that node from the editor. Only the endpoint nodes may not be deleted.

Below the color map editor (or transfer function editor, for volume rendering) is a set of three text boxes for changing the scalar value, opacity, and scale associated with a given node. Only the scalar value is associated with surface rendering. The scalar values at the endpoints may only be changed if the **Automatically Rescale to Fit Data Range** check box (discussed later in this section) is unmarked. When volume rendering, you may also specify the opacity and scale per node in the editor. In volume rendering, the opacity is accumulated as you step through the volume being rendered. The **Scale** value determines the unit distance over which the opacity is accumulated.

The controls below the set of three text boxes specify the color space and any color map preset you wish to save or use. The color spaces available are **RGB** (red, green, blue), **HSV** (hue, saturation, value), **Wrapped HSV**, and **CIELAB** (a more perceptually linear color space). The color space determines how the colors are interpolated between specified values; the colors at the color map (or transfer function) editor nodes will remain the same regardless of the color space chosen. If wrapped HSV is used, the interpolation will use the shortest path in hue, even going through the value hue = 0. For non-wrapped HSV, the hue interpolation will not pass through 0. (A hue of zero sets the color to red.)

In addition to choosing the color space and modifying the color map / transfer function nodes, you may also create and load preset color scales. (When volume rendering, only the color map is stored; the scalar-to-opacity mapping is not.) To store your current settings as a preset, click the **Save** button. In the dialog box that appears, you may enter a name for your new preset. By default, the scalar values from the data array being used are stored in the preset. If you wish these values to be normalized between 0 and 1, press the **Normalize** button.

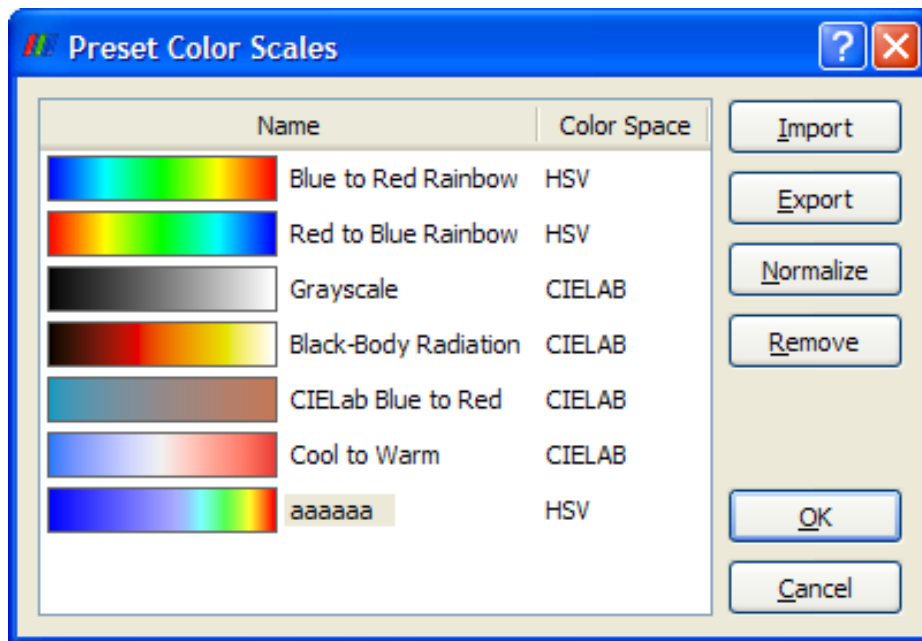


Figure 2. Dialog for choosing color scale presets.

Any presets you save, in addition to the default ones provided by ParaView, are available by pressing the **Choose Preset** button, causing the dialog shown below to be displayed. Selecting a preset and clicking **OK** causes the current color map to be set to the chosen preset. Any user-defined presets may be normalized (as discussed above) or removed from the list of presets entirely using the **Normalize** and **Remove** buttons, respectively. (The default presets are already normalized and may not be removed from the application.)

Any of the color scale presets may be exported to a file using the **Export** button in the above dialog. The resulting file(s) may then be copied to another computer for use with ParaView on a different machine. In order to load presets that are stored in such files, press the **Import** button on the above dialog, and navigate to the desired color preset file.

If the current data set is colored by an array of vectors, the **Component** menu will be enabled. It determines whether the data is colored by a single vector component (**X**, **Y**, or **Z**) or by the vector's **Magnitude** (the default). If the data is colored by a single-component (scalar) array, then the **Component** menu is disabled.

If **Use Logarithmic Scale** is checked, then instead of the scalar values in the data array being used directly to determine the colors, the base-10 logarithm of the data array values is computed, and the resulting value is used for extracting a color from the color map. If the data array contains values for which a logarithm would produce invalid results (i.e., any values less than or equal to 0), the range for the color map is changed to [0, 10] so that the logarithm produces valid results.

By default, any data attribute that has been used to color a data set currently loaded in ParaView whose name and number of components match that of the array selected in the **Color by** menu contributes to the range of the color map. To change this behavior, first uncheck the **Automatically Rescale to Fit Data Range** check box. (This ensures that the range of the color map is not reset when the range of the data attribute changes.) The minimum and maximum values of the color map can be overridden by pressing the **Rescale Range** button, entering different **Minimum** and **Maximum** values in the dialog that appears, and pressing **Rescale**; this rescales all the nodes in the color map so that the scalar values lie at the same normalized positions. Alternatively, you may modify the scalar values of any node (including the endpoints if **Automatically Rescale to Fit Data Range** is off) by clicking a node to highlight it and typing a new value in the **Scalar Value** entry box. By changing the minimum and maximum color map values, it is possible to manually specify what range of data values the color map will cover. Pressing the **Rescale to Data Range** button on the **Color Scale** tab of the **Color Scale Editor** sets the range to cover only the current data set.

If **Use Discrete Colors** is checked, the **Resolution** slider at the bottom of the dialog specifies the number of colors to in the color map. The scale ranges from 2 to 256 (the default). The fewer the number of colors, the larger the range each color covers. This is useful if the data attribute has a small number of distinct values or if larger ranges of the array values should be mapped to the same color.

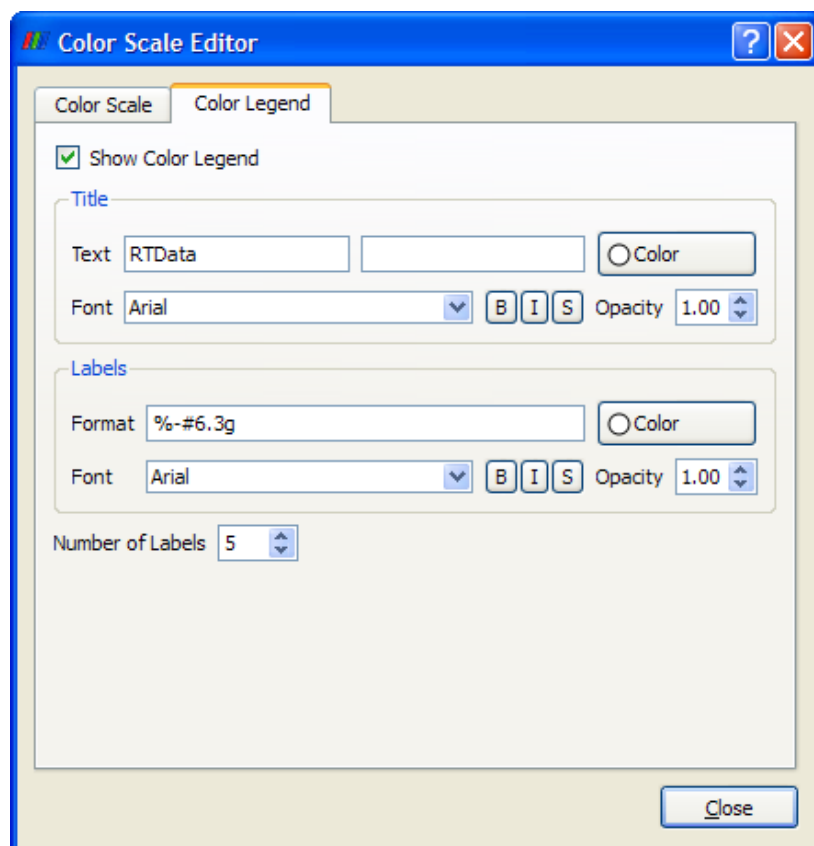


Figure 3. Scalar Bar controls



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# Modifying Data

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## Rationale

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### Manipulating Data

In the course of either searching for information within data or in preparing images for publication that help to explain data, it is often necessary to process the raw data in various ways. Examples include slicing into the data to make the interior visible, extracting regions that have particular qualities, and computing statistical measurements from the data. All of these operations involving taking in some original data and using it to compute some derived quantities. This chapter explains how one controls the data processing portion of ParaView's visualization pipeline to go about doing such analyses.

A filter is the basic tool that one uses to manipulate data. If data is a noun, then a filter is the verb that operates on the data. Filters operate by ingesting some data, processing it and producing some other data. In the abstract sense a data reader is a filter too because it ingests data from the file system. ParaView creates filters then when you open data files, and when you instantiate new filters from the Filters menu. The set of filters you create becomes your visualization pipeline, and that pipeline is shown in ParaView's Pipeline Browser.

## Filter Parameters

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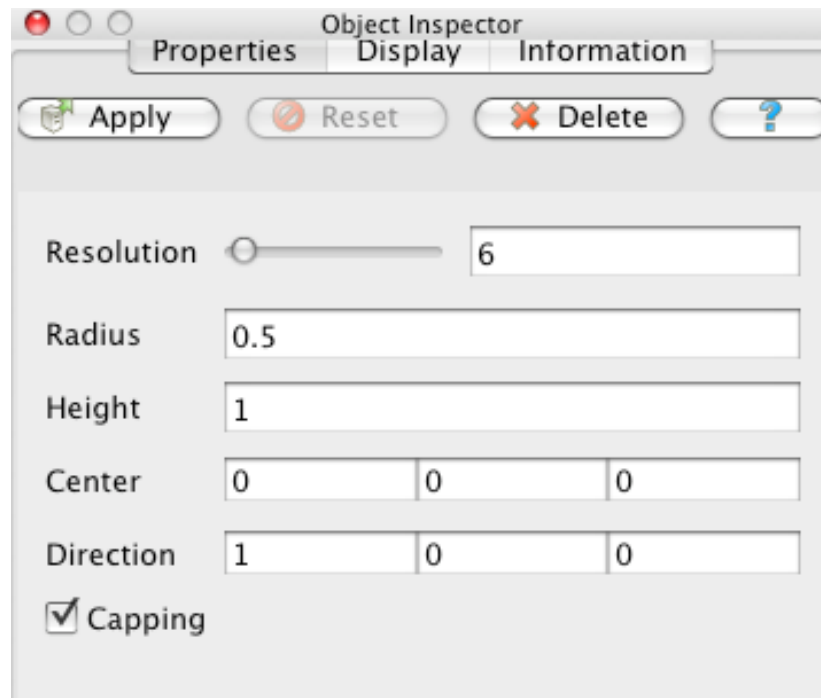
### Filter Parameters

Each time a data set is opened from a file, a source is selected, a filter is applied, or an existing reader, source, or filter (hereafter simply referred to as filter) is selected in the **Pipeline Browser**, ParaView updates the **Object Inspector** for the corresponding output data set. The **Object Inspector** has three tabs. In this chapter we are primarily concerned with the **Properties** tab. The **Display** tab gives you control over the visual characteristics of the data produced by the filter as displayed in the active view and the **Information** tab presents meta-information about the data produced by the filter.

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## Properties

From the *Properties*' tab you modify the parameters of the filter, fine tuning what it produces from its input (if any). A filter that extracts an isocontour will have a control with which to set the isovalue (or isovalues) to extract at for example. The specific controls and information provided on the tab then are specific to the particular `vtkAlgorithm` you are working with, but all filters have at least the **Apply**, **Reset**, **Delete** and ? (help) controls.



**Figure 1. Sample Properties tab for a Cone source**

The help button brings opens up the documentation for the filter in ParaView's help system in which the input restrictions to the filter, output type generated by the filter, and descriptions of each parameter are listed. The same information is repeated in the Appendices 1,2 of this book.

The **Delete** button removes this filter from the pipeline. The delete button is only enabled when there are no filters further down the pipeline that depend on this filter's output. You have to either use the Pipeline Browser and Object inspector in conjunction to delete the dependent parts of the pipeline or use **Delete All** from the **Edit** menu.

When a reader, source, or filter is first selected, the associated data set is not immediately created. By default (unless you turn on **Auto-Accept** in ParaView's settings) the filter will not run until you hit the **Apply** button. When you do press Apply, ParaView sends the values shown on the **Properties** tab to the data processing engine and then the pipeline is executed. This delayed commit behavior is important when working with large data, for which any given action might take a long time to finish.

Until you press **Apply** and any other time that the values shown on the GUI do not agree with what was last sent to the server, the the Apply button will be highlighted (in blue or green depending on your operating system). In this state the **Reset** button is also enabled. Pressing that returns the GUI to the last committed state, which gives you an easy way to cancel mistakes you've made before they happen.

The specific parameter control widgets vary from filter to filter and sometimes vary depending on the exact input to the filter. Some filters have no parameters at all and others have many. Many readers present the list and type of arrays in the file, an allow you to pick some or all of them as you need. In all cases the widgets shown on the **Properties** tab give you control over exactly what the filter does. If you are unsure of what they do, remember to hit the ? button to see the documentation for that filter.

Note that ParaView attempts to provide reasonable default settings for the parameter settings and to some extent guards against invalid entries. A numeric entry box will not let you type in non-numerical values for example.

Sliders and spin boxes typically have minimum and maximum limits built in. In some cases though you may want to ignore these default limits. Whenever there is a numeric entry beside the widget, you are able to manually type in any number you need to.

Some filter parameters are best manipulated directly in the 3D View window with the mouse. Take the Slice filter for example. Slice extracts slices from the data that lie on a set of parallel planes oriented in space. This type of world space interactive control is what 3D Widgets are for. The textual controls in the Properties Tab and the displayed state of the 3D widgets are always linked so that changing either updates the other. You can of course use the **Reset** button to revert changes you make from either place.

## The Pipeline

### Managing the Pipeline

Data manipulation in ParaView is fairly unique because of the underlying pipeline architecture that it inherits from VTK. Each filter takes in some data and produces something new from it. Filters do not directly modify the data that is given to them and instead copy unmodified data through via reference (so as to conserve memory) and augment that with newly generated or changed data. The fact that input data is never altered in place means that unlike many visualization tools, you can apply several filtering operations in different combinations to your data during a single ParaView session. You see the result of each filter as it is applied, which helps to guide your data exploration work, and can easily display any or all intermediate pipeline outputs simultaneously.

The **Pipeline Browser** depicts ParaView's current visualization pipeline and allows you to easily navigate to the various readers, sources, and filters it contains. Connecting an initial data set loaded from a file or created from a ParaView source to a filter creates a two filter long visualization pipeline. The initial data set read in becomes the input to the filter, and if needed the output of that filter can be used as the input to a subsequent filter, etc.

For example, suppose you create a sphere in ParaView by selecting **Sphere** from the **Sources** menu. In this example, the sphere is the initial data set. Next create a **Shrink** filter from the **Alphabetical** submenu of the **Filters** menu. Because the sphere source was the active filter when the shrink filter was created, the shrink filter operates on the sphere source's output. Optionally, use the **Properties** tab of the **Object Inspector** to set initial parameters of the shrink filter and then hit **Apply**. Next create an **Elevation** filter to filter the output of the shrink filter and hit **Apply** again. You have just created a simple three element linear pipeline in ParaView. You will now see the following in the **Pipeline Browser**.

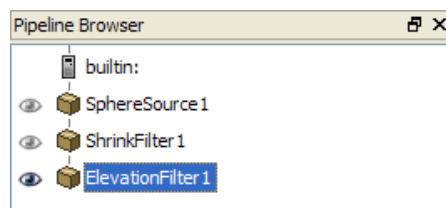
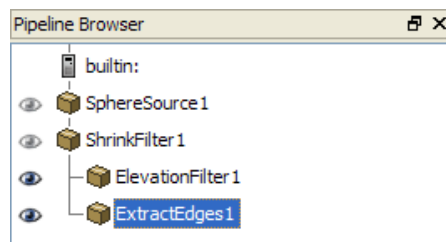


Figure 2. Linear Pipeline

Within the **Pipeline Browser**, to the left of each entry is an **eye** icon indicating whether that data set is currently visible. If there is no **eye** icon, it means that the data produced by that filter is not compatible with the currently active View window. Otherwise, a dark **eye** icon indicates that the data set is visible; and when a data set is viewable but currently invisible, its icon is drawn in light gray. Clicking on the **eye** icon toggles the visibility of the corresponding data set. In the above example, all three filters are potentially visible, but only the **ElevationFilter** is actually being displayed. The **ElevationFilter** is also highlighted in blue, indicating that it is the currently "active" Filter. Since it is the "active" filter, the **Object Inspector** reflects its content and the next filter created will use it as the input.

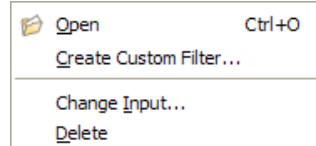
You can always change parameters of any filter in the pipeline after it has been applied. Left-click on the filter in the **Pipeline Browser** to make it the "active" one. The **Properties**, **Display**, and **Information** tabs always reflect the "active" filter. When you make changes in the **Properties** tab and apply your changes, all filters beyond the changed one are automatically updated. Double-clicking the name of one of the filters causes the name to become editable, enabling you to change it to something more meaningful than the default chosen by ParaView.

By default each filter you add to the pipeline becomes the active filter, which is handy when making linear pipelines. Branching pipelines are also very useful. The simplest way to make one is to click on some other, further upstream filter in the pipeline before you create a new filter. For example Select ShrinkFilter1 in the **Pipeline Browser** then apply **Extract Edges** from the **Alphabetical** submenu of the **Filters** menu. Now the output of the shrink filter is being used as the input to both the elevation and extract edges filters. You will see the following in the Pipeline Browser.



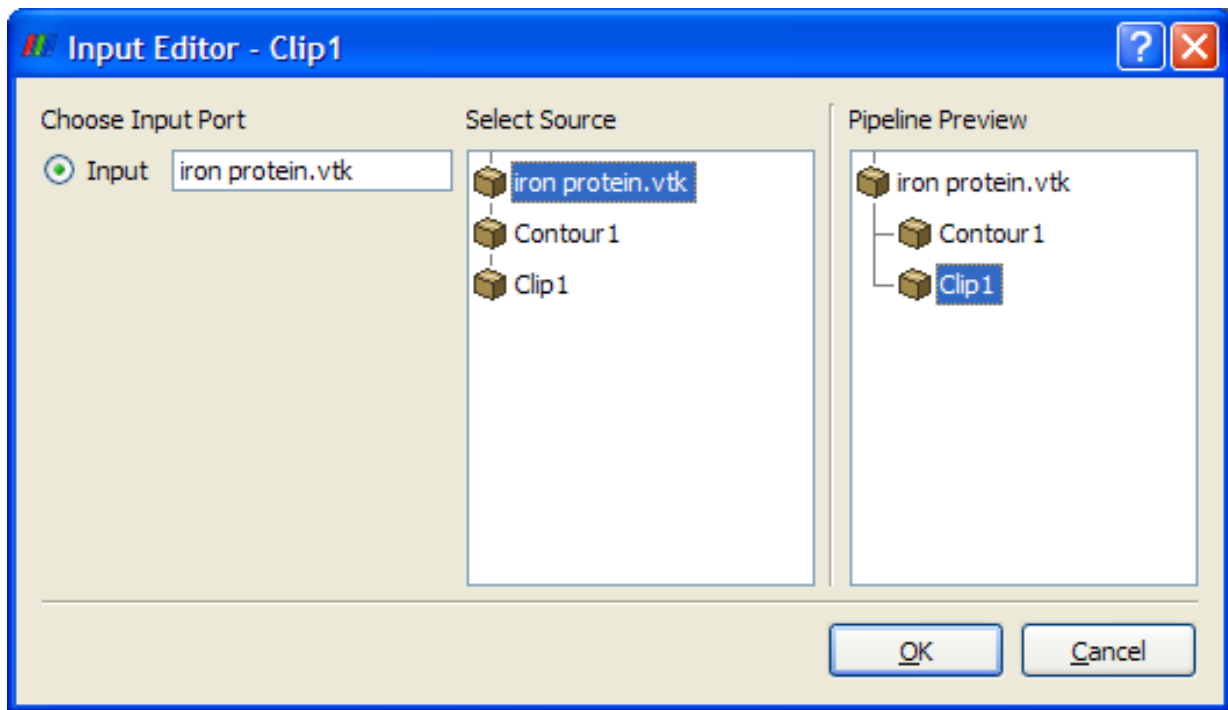
**Figure 3. Branching Pipeline**

Right-clicking a filter in the *Pipeline Browser* displays a context menu from which you can do several things. For reader modules you can use this to load a new data file. All modules can be saved (the filter and the parameters you've set on it) as a Custom Filter (see the last section of this chapter), or delete it (if it is at the end of the visualization pipeline). For filter modules you can also use this menu to change the input to the filter, and thus rearrange the visualization pipeline.



**Figure 4. Context menu in Pipeline Browser**

To rearrange the pipeline select **Change Input** from the context menu. That will bring up the **Input Editor** dialog box shown in **Figure 5**. The name of the dialog box reflects the filter that you are moving. The middle **Select Source** pane shows the pipeline as it stands currently and use this pane to select a filter to move the chosen filter on to. This pane only allows you to choose compatible filters. I.e, ones that produce data that can be ingested by the filter you are moving. It also does not allow you to create loops in the pipeline. Left-click to make your choice from the allowed set of filter and then the rightmost **Pipeline Preview** pane will show what the pipeline will look like once you commit your change. Click **OK** to do so or **Cancel** to abort the move.



**Figure 5. Input Editor dialog shown while moving an elevation filter to operate directly on a reader.**

Some filters require more than one input. (See the discussion of merging pipelines below). For those the leftmost **input port** pane of the dialog produces shows more than one port. Use that together with the **Select Source** pane to specify each input in turn.

Conversely, some filters produce more than one output. Thus another way to make a branching pipeline is simply to open a reader for instance that produces multiple distinct data sets. An example of this is the SLAC reader that produces both a polygonal output and a structured data field output. Whenever you have a branching pipeline keep in mind that it is important to select the proper branch on which to extend the pipeline. For example, if you want to apply a filter like the Extract Subset filter, which operates only on structured data, while SLAC reader's polygonal output is the currently active filter, until you click on the reader's structured data output and make that the currently active one.

Some filters that produce multiple data sets do so in a different way. Instead of producing several fundamentally distinct data sets, they produce a single composite data set which contains many sub data sets. See the Understanding Data chapter for an explanation of composite data sets. With composite data it is usually best to treat the entire group as one entity and you do not need to do anything in particular to do so. Sometimes though, you want to operate on a particular set of sub data sets. To do that apply the **Extract Block** filter. This filter allows you to pick the desired sub data set(s). Next apply the filter you are actually interested in to the extract filter's output. An alternative is to hit 'B' to use '[ParaView/Users Guide/Selection/Block Selection]' in a 3D View and then use the **Extract Selection** filter.

Pipelines merge as well, whenever they contain filters that take in more than one input to produce their own output (or outputs). There are in fact two different varieties of merging filters. The **Append Geometry** and **Group Data Sets** filters are examples of the first kind. These filters take in any number of fundamentally similar data sets. Append for example takes in one or more polygonal datasets and combines them into a single large polygonal data set. Group takes in a set of any type of datasets and produces a single composite dataset. To use this type of merging filter, select more than one filter within the Pipeline browser by left clicking to select the first input and then shift-left clicking to select the rest. Now create the merging filter from the **Filters** menu as usual. The pipeline in this case will look like the one in the following figure.

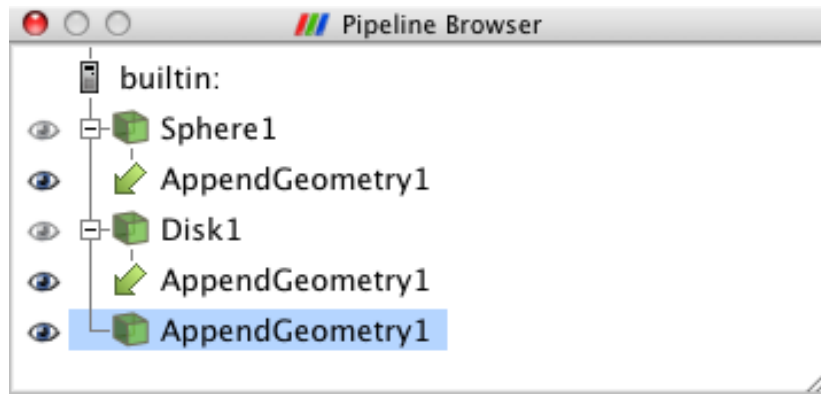


Figure 7. merging pipeline

Other filters take in more than one, fundamentally different data sets. An example is the **Resample with Dataset** filter which takes in one data set (the **Input**) that represents a field in space to sample values from and another data set (the **Source**) to use as a set of locations in space to sample the values at. Begin this type of merge by choosing either of the two inputs as the active filter and then creating the merging filter from the **Filters** menu. A modified version of the **Change Input** dialog shown in **Figure 5** results (this one that lacks a **Pipeline Preview** pane). Click on either of the ports listed in the **Available Inputs** pane and specify an input for it from the **Select Input** pane. Then switch to the other input in the **Available Inputs** port and choose the other input on the **Select Input** pane. When you are satisfied with your choices click **OK** on the dialog and then **Apply** on the **Pipeline Browser** to create the merging filter.

## Filter Categories

### Available Filters

There are a great many filters available in ParaView (1) (and even more in VTK), and because ParaView has a modular architecture, it is routine for people to add additional ones (2). Some filters have obscure purposes and are rarely used, but others are more general purpose and used very frequently. These most common filters are found easily on the **Common (View -> Toolbars)** toolbar.



Figure 10. Common Filters Toolbar

These filters include:

- **Calculator** - Evaluates a user-defined expression on a per-point or per-cell basis (3)
- **Contour** - Extracts the points, curves, or surfaces where a scalar field is equal to a user-defined value. This surface is often also called an isosurface.
- **Clip** - Intersects the geometry with a half space. The effect is to remove all the geometry on one side of a user-defined plane.
- **Slice** - Intersects the geometry with a plane. The effect is similar to clipping except that all that remains is the geometry where the plane is located.
- **Threshold** - Extracts cells that lie within a specified range of a scalar field.
- **Extract Subset** - Extracts a subset of a grid by defining either a volume of interest or a sampling rate.
- **Glyph** - Places a glyph, a simple shape, on each point in a mesh. The glyphs may be oriented by a vector and scaled by a vector or scalar.
- **Stream Tracer** - Seeds a vector field with points and then traces those seed points through the (steady state) vector field.

- **Warp** - Displaces each point in a mesh by a given vector field.
- **Group Datasets** - Combines the output of several pipeline objects into a single multi block data set.
- **Group Extract Level** - Extract one or more items from a multi block data set.

These eleven filters are a small sampling of what is available in ParaView.

In the Alphabetical submenu of the Filters menu you will find all of the filters that are useable in your copy of ParaView. Currently there are more than one hundred of them, so to make them easier to find the Filters menu is organized into submenus. These submenus are organized as follows.

- **Recent** - The filters you've used recently.
- **Common** - The common filters. This is the same set of filters on the common filters toolbar.
- **Cosmology** - This contains filters developed at LANL for cosmology research.
- **Data Analysis** - The filters designed to retrieve quantitative values from the data. These filters compute data on the mesh, extract elements from the mesh, or plot data.
- **Statistics** - This contains filters that provide descriptive statistics of data, primarily in tabular form.
- **Temporal** - Filters that analyze or modify data that changes over time.

All filters can work on data that changes over time because they are re-executed at each time step. Filters in this category have the additional capability to inspect and make use of or even modify the temporal dimension.

- **Alphabetical** - Many filters do not fit into the above categories so all filters can be found here.

| Filter                  | Window                | Help                    |
|-------------------------|-----------------------|-------------------------|
| All to N                | Extract Parts         | Part Id Scalars         |
| Append Attributes       | Extract Surface       | Pick                    |
| Append Datasets         | Feature Edges         | Point Data to Cell Data |
| Append Geometry         | Generic Outline       | Probe                   |
| Balance                 | Generic Stream Tracer | Process Id Scalars      |
| Calculator              | GenericClip           | Quadric Clustering      |
| Cell Centers            | GenericContour        | Random Vectors          |
| Cell Data to Point Data | GenericCut            | Reflection              |
| Clean                   | GenericGeometryFilter | Ribbon                  |
| Clean to Grid           | Glyph                 | Rotational Extrusion    |
| Clip                    | Gradient              | Shrink                  |
| Connectivity            | Gradient Magnitude    | Smooth                  |
| Contour                 | Group Datasets        | Stream Tracer           |
| Crop                    | Group Parts           | Subdivide               |
| Curvature               | Integrate Attributes  | Surface Flow            |
| Cut                     | Level Scalars         | Surface Vectors         |
| D3                      | Linear Extrusion      | Tessellate              |
| Decimate                | Loop Subdivision      | Tessellator             |
| Delaunay2D              | Mask Points           | Tetrahedralize          |
| Elevation               | Median                | Threshold               |
| Extract CTH Parts       | Mesh Quality          | Transform               |
| Extract Datasets        | Normals generation    | Triangle Strips         |
| Extract Edges           | Outline               | Triangulate             |
| Extract Grid            | Outline (curvilinear) | Tube                    |
| Extract Group           | Outline Corners       | Warp (scalar)           |
|                         |                       | Warp (vector)           |

**Figure 11. A portion of the Alphabetical submenu of the Filters menu.**

Searching through these lists of filters, particularly the full alphabetical list, can be cumbersome. To speed up the selection of filters, you should use the quick launch dialog. Choose the first item from the filters menu, or alternatively press either **CTRL** and **SPACE BAR** (Windows or Linux) or **ALT** and **SPACE BAR** (on Macintosh) together to bring up the **Quick Launch** dialog. As you type in words or word fragments the dialog lists the filters whose names contain them. Use the up and down arrow key to select from among them and hit **ENTER** to create the filter.

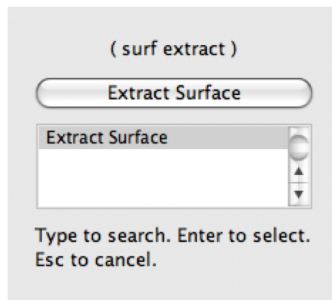


Figure 12. Quick Launch

### Why can't I apply the filter I want?

Note that many of the filters in the menu will be grayed out and not selectable at any given time. That is because any given filter may only operate on particular types of data. For example the Extract Subset will only operate on structured data sets so it is only enabled when the module you are building on top of produces image data, rectilinear grid data, or structured grid data. Likewise the contour filter requires scalar data and can not operate directly on data sets that have only vectors. The input restrictions for all filters are listed in the Appendix and help menus.

When the filter you want is not available you should look for a similar filter which will accept your data or apply an intermediate filter which transforms your data into the required format. Since ParaView 3.10 you can also ask ParaView to try to do the conversion for you automatically by clicking "Auto convert Properties" in the application settings.

### What does that filter do?

A description of what each filter does, what input data types it accepts and what output data types it produces can be found in the [ParaViewUsersGuide/List\_of\_filters/Appendix] and help menus. For a more complete understanding, remember that most ParaView filters are simply VTK algorithms, each of which is documented online in the VTK (<http://www.vtk.org/doc/release/5.6/html/classes.html>) and ParaView (<http://www.paraview.org/ParaView3/Doc/Nightly/html/classes.html>) Doxygen wiki pages.

When one is exploring a given data set, one does not want to have to hunt through the detailed descriptions of all of the filters in order to find the one filter that is needed at any given moment. It is useful then to be aware of the general high-level taxonomy of the different operations that the filters can be logically grouped into.

These are:

- Attribute Manipulation : which manipulate the field aligned, point aligned and cell aligned data values and in general derive new aligned quantities.

Including **Curvature**, **Elevation**, **Generate Ids**, **Generate Surface Normals**, **Gradient**, **Mesh Quality**, **Principal Component Analysis**, and **Random Vectors**.

- Geometric Manipulation : which essentially operate on or manipulate the shape of the data in a spatial context.

Including **Reflect**, **Transform**, and **Warp**

- Topological operations : which manipulate the connected structure of the data set itself, usually creating or destroying cells, for instance to reduce the data sets memory size while leaving it in the same place in space.

Including **Cell Centers**, **Clean**, **Decimate**, **Extract Surface**, **Quadric Clustering**, **Shrink**, **Smooth**, and **Tetrahedralize**.

- Sampling : which compute new data sets that represent some essential features from the data sets that they take as input.

Including **Clip**, **Extract Subset**, **Extract Selection**, **Glyph**, **Streamline**, **Probe**, **Plot**, **Histogram**, and **Slice**



- **Data Type Conversion** : which convert between the various VTK data structures VTK Data Model and join or split entire data structures

Including **Append DataSets**, **Append Geometry**, **Extract Blocks**, **Extract AMR Blocks**, and **Group DataSets**

- **White Box Filters** : which do arbitrary processing as specified at runtime by you the user

Including the **Calculator** and **Python Programmable** filters.

## Best Practices

---

### 1.4.1 Avoiding Data Explosion

The pipeline model that ParaView presents is very convenient for exploratory visualization. The loose coupling between components provides a very flexible framework for building unique visualizations, and the pipeline structure allows you to tweak parameters quickly and easily.

The downside of this coupling is that it can have a larger memory footprint. Each stage of this pipeline maintains its own copy of the data. Whenever possible, ParaView performs shallow copies of the data so that different stages of the pipeline point to the same block of data in memory. However, any filter that creates new data or changes the values or topology of the data must allocate new memory for the result. If ParaView is filtering a very large mesh, inappropriate use of filters can quickly deplete all available memory. Therefore, when visualizing large data sets, it is important to understand the memory requirements of filters.

Please keep in mind that the following advice is intended only for when dealing with very large amounts of data and the remaining available memory is low. When you are not in danger of running out of memory, ignore all of the following advice.

When dealing with structured data, it is absolutely important to know what filters will change the data to unstructured. Unstructured data has a much higher memory footprint, per cell, than structured data because the topology must be explicitly written out. There are many filters in ParaView that will change the topology in some way, and these filters will write out the data as an unstructured grid, because that is the only data set that will handle any type of topology that is generated. The following list of filters will write out a new unstructured topology in its output that is roughly equivalent to the input. These filters should never be used with structured data and should be used with caution on unstructured data.

- **Append Datasets**
  - **Append Geometry**
  - **Clean**
  - **Clean to Grid**
  - **Connectivity**
  - **D3**
  - **Delaunay 2D/3D**
  - **Extract Edges**
  - **Linear Extrusion**
  - **Loop Subdivision**
  - **Reflect**
  - **Rotational Extrusion**
  - **Shrink**
  - **Smooth**
  - **Subdivide**
  - **Tessellate**
  - **Tetrahedralize**
-

- **Triangle Strips**
- **Triangulate**

Technically, the **Ribbon** and **Tube** filters should fall into this list. However, as they only work on 1D cells in poly data, the input data is usually small and of little concern.

This similar set of filters also output unstructured grids, but they also tend to reduce some of this data. Be aware though that this data reduction is often smaller than the overhead of converting to unstructured data. Also note that the reduction is often not well balanced. It is possible (often likely) that a single process may not lose any cells. Thus, these filters should be used with caution on unstructured data and extreme caution on structured data.

- **Clip**
- **Decimate**
- **Extract Cells by Region**
- **Extract Selection**
- **Quadric Clustering**
- **Threshold**

Similar to the items in the preceding list, **Extract Subset** performs data reduction on a structured data set, but also outputs a structured data set. So the warning about creating new data still applies, but you do not have to worry about converting to an unstructured grid.

This next set of filters also outputs unstructured data, but it also performs a reduction on the dimension of the data (for example 3D to 2D), which results in a much smaller output. Thus, these filters are usually safe to use with unstructured data and require only mild caution with structured data.

- **Cell Centers**
- **Contour**
- **Extract CTH Fragments**
- **Extract CTH Parts**
- **Extract Surface**
- **Feature Edges**
- **Mask Points**
- **Outline (curvilinear)**
- **Slice**
- **Stream Tracer**

These filters do not change the connectivity of the data at all. Instead, they only add field arrays to the data. All the existing data is shallow copied. These filters are usually safe to use on all data.

- **Block Scalars**
  - **Calculator**
  - **Cell Data to Point Data**
  - **Curvature**
  - **Elevation**
  - **Generate Surface Normals**
  - **Gradient**
  - **Level Scalars**
  - **Median**
  - **Mesh Quality**
  - **Octree Depth Limit**
  - **Octree Depth Scalars**
  - **Point Data to Cell Data**
  - **Process Id Scalars**
-

- **Random Vectors**
- **Resample with dataset**
- **Surface Flow**
- **Surface Vectors**
- **Texture Map to...**
- **Transform**
- **Warp (scalar)**
- **Warp (vector)**

This final set of filters are those that either add no data to the output (all data of consequence is shallow copied) or the data they add is generally independent of the size of the input. These are almost always safe to add under any circumstances (although they may take a lot of time).

- **Annotate Time**
- **Append Attributes**
- **Extract Block**
- **Extract Datasets**
- **Extract Level**
- **Glyph**
- **Group Datasets**
- **Histogram**
- **Integrate Variables**
- **Normal Glyphs**
- **Outline**
- **Outline Corners**
- **Plot Global Variables Over Time**
- **Plot Over Line**
- **Plot Selection Over Time**
- **Probe Location**
- **Temporal Shift Scale**
- **Temporal Snap-to-Time-Steps**
- **Temporal Statistics**

There are a few special case filters that do not fit well into any of the previous classes. Some of the filters, currently **Temporal Interpolator** and **Particle Tracer**, perform calculations based on how data changes over time. Thus, these filters may need to load data for two or more instances of time, which can double or more the amount of data needed in memory. The **Temporal Cache** filter will also hold data for multiple instances of time. Also keep in mind that some of the temporal filters such as the temporal statistics and the filters that plot over time may need to iteratively load all data from disk. Thus, it may take an impractically long amount of time even if does not require any extra memory.

The **Programmable Filter** is also a special case that is impossible to classify. Since this filter does whatever it is programmed to do, it can fall into any one of these categories.

### 1.4.2 Culling Data

When dealing with large data, it is clearly best to cull out data whenever possible, and the earlier the better. Most large data starts as 3D geometry and the desired geometry is often a surface. As surfaces usually have a much smaller memory footprint than the volumes that they are derived from, it is best to convert to a surface soon. Once you do that, you can apply other filters in relative safety.

A very common visualization operation is to extract isosurfaces from a volume using the contour filter. The **Contour** filter usually outputs geometry much smaller than its input. Thus, the **Contour** filter should be applied early if it is to

be used at all. Be careful when setting up the parameters to the **Contour** filter because it still is possible for it to generate a lot of data. This obviously can happen if you specify many isosurface values. High frequencies such as noise around an isosurface value can also cause a large, irregular surface to form.

Another way to peer inside of a volume is to perform a **Slice** on it. The **Slice** filter will intersect a volume with a plane and allow you to see the data in the volume where the plane intersects. If you know the relative location of an interesting feature in your large data set, slicing is a good way to view it.

If you have little a-priori knowledge of your data and would like to explore the data without paying the memory and processing time for the full data set, you can use the **Extract Subset** filter to subsample the data. The subsampled data can be dramatically smaller than the original data and should still be well load balanced. Of course, be aware that you may miss small features if the subsampling steps over them and that once you find a feature you should go back and visualize it with the full data set.

There are also several features that can pull out a subset of a volume: **Clip**, **Threshold**, **Extract Selection**, and **Extract Subset** can all extract cells based on some criterion. Be aware, however, that the extracted cells are almost never well balanced; expect some processes to have no cells removed. Also, all of these filters with the exception of **Extract Subset** will convert structured data types to unstructured grids. Therefore, they should not be used unless the extracted cells are of at least an order of magnitude less than the source data.

When possible, replace the use of a filter that extracts 3D data with one that will extract 2D surfaces. For example, if you are interested in a plane through the data, use the **Slice** filter rather than the **Clip** filter. If you are interested in knowing the location of a region of cells containing a particular range of values, consider using the **Contour** filter to generate surfaces at the ends of the range rather than extract all of the cells with the **Threshold** filter. Be aware that substituting filters can have an effect on downstream filters. For example, running the **Histogram** filter after **Threshold** will have an entirely different effect than running it after the roughly equivalent **Contour** filter.

## Custom Filters aka Macro Filters

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### Macros (aka Custom Filters)

It often happens that once you figure out how to do some specific data processing task, you want to repeat it often. You may for example want to reuse particular filters with specific settings (for example complicated calculator or programmable filter expressions) or even entire pipeline sections consisting on new data sets without having to manually enter the parameters each time.

You can do this via the clever use of state files or more conveniently python scripts and python macros 1 . Saving, editing and reusing state files gives you the ability to recreate entire ParaView sessions. This does not give you fine enough control for small, repeatedly reused tasks. Python Tracing does give you fine grained control, but this assumes that you have python enabled in your copy of ParaView (which is usually but not always the case) and that you remember to turn on Trace recording before you did whatever it was that you want to play back. Both techniques largely require that you think like a programmer when you initially create and setup the scripts. Another alternative is to use ParaView's Custom Filters which let you create reusable meta-filters strictly within the GUI.

A Custom Filter is a black box filter that encapsulates one or more filters in a sub-pipeline and exposes only those parameters from that sub-pipeline that the Custom Filter creator chose to make available. For example, if you capture a ten element pipeline in your Custom Filter where each filter happened to have eight parameters, you could choose to expose anywhere from zero to eighty parameters in your Custom Filter's **Properties** tab.

---

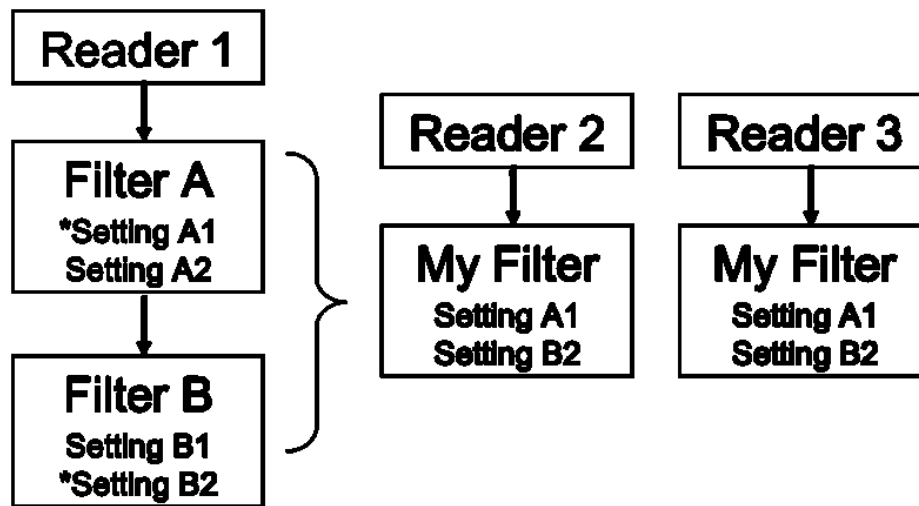
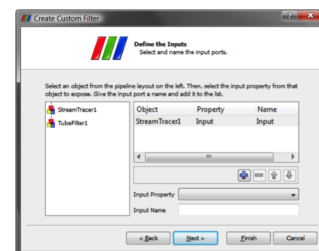


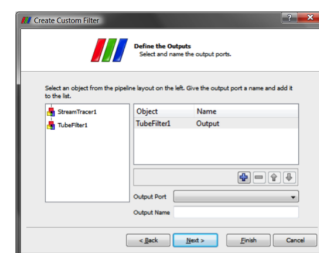
Figure 13: Custom Filter concept

Once you have set up some pipeline that performs the data processing that you want to reuse, the process of creating a Custom Filter consists of three steps. First select one or more filters from the **Pipeline Browser** using the mouse. Next from the **Tools** menu select **Create Custom Filter**. From that dialog choose the filter in your chosen sub-pipeline who's input is representative of where you want data to enter into your Custom Filter. This is usually the topmost filter. If you are creating a multi-input filter, click the **+** button to add additional inputs and configure them in the same way. Clicking **Next** brings you to a similar dialog in which you choose the outputs of your Custom Filter. Click **Next** again to get to the last dialog. On this one you specify which parameters of the internal filters that you want to expose to the eventual user of the custom filter. You can optionally give each parameter a descriptive label here as well. The three dialogs are shown below.

Step 1: configure one or more inputs to your new filter.



Step 2: configure one or more outputs of your new filter.



Step 3: identify and name the controls you want to expose of your new filter.

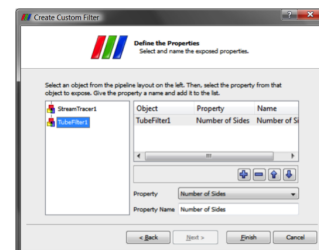


Figure 9: Creating a Custom Filter

Once you create a Custom Filter it is added to the **Alphabetical** sub menu of the **Filters** menu. It is automatically saved in ParaView's settings, so the next time you start ParaView on the same machine you can use it just like any of

the other filters that came with your copy of ParaView. Custom Filters are treated no differently than other filters in ParaView and are saveable and restorable in state files and python scripts. If you find that you no longer need some Custom Filter and want to get rid of it, use the **Manage Custom Filters** dialog box under the **Tools** menu to remove it.

If on the other hand you find that a Custom Filter is very useful, you may instead want to give it to a colleague. On that same dialog are controls for exporting and importing Custom Filters. When you save a Custom Filter you are prompted for a location and filename to save the filter in. The file format is a simple XML text file that you can simply email or otherwise deliver.

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# Quantative Analysis

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## Drilling Down

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ParaView 2 was almost entirely a qualitative analysis tool. It was very good at drawing pictures of large scientific data sets so that you could view the data and tell if it looked "right", but it was not easy to use for finding hard quantitative information about the data and verifying that that was the case. The recommended use was to use ParaView to visualize, interact with and subset your data and then export the result into a format that could be imported by a different tool. A major goal of ParaView 3 has been to add quantitative analysis capabilities to turn it into a convenient and comprehensive tool in which you can visualize, interact with and drill all the way down into the data.

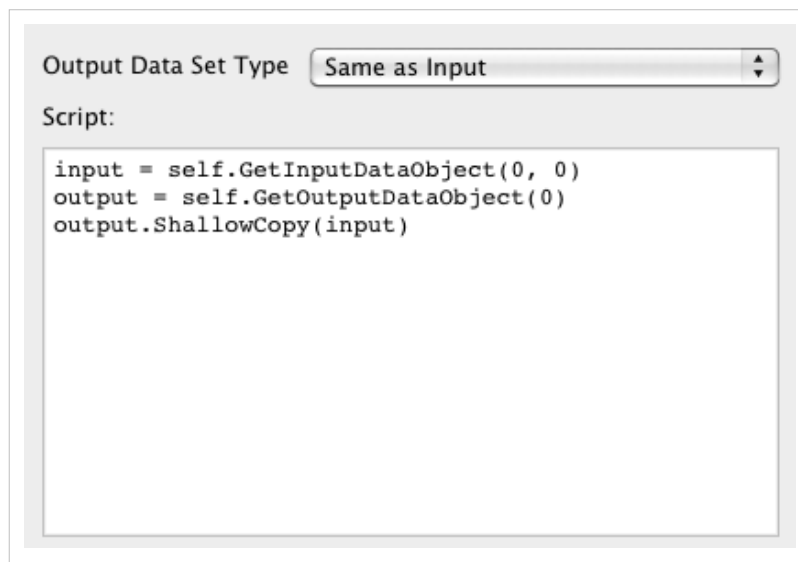
These capabilities vary from semi-qualitative ones such as the **Ruler** source and **Cube Axis** representation (see [Users\\_Guide\\_Annotation](#)) to Selection which allows you to define and extract arbitrary subsets of the data based, to the spreadsheet view which presents the data in textual form. Taken together with features like the statistical analysis filters [ParaView/Users\\_Guide/List\\_of\\_filters#Descriptive\\_Statistics](#), calculator filters, 2D plot and chart views and programmable filters (which give you the ability to run arbitrary code on the server and have access to every data point) these give you the ability to inspect the data at from the highest level view all the way down to the hard numbers.

In this chapter we describe the various tools that ParaView gives you to support quantitative analysis.

## Python Programmable Filter

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### Introduction



The Programmable Filter is a ParaView filter that processes one or more input datasets based on a Python script provided by the user. The parameters of the filter include the output data type, the script and a toggle that controls whether the input arrays are copied to the output. In this document, we introduce the use of the Programmable Filter and give a summary of the API available to the user.

Note that the Programmable Filter depends on Python. All ParaView binaries distributed by Kitware are

built with Python enabled. If you have built ParaView yourself, you have to make sure that `PARAVIEW_ENABLE_PYTHON` is turned on when configuring the ParaView build.

---

Since the entire VTK API as well as any module that can be imported through Python is available through this filter, we can only skim the surface of what can be accomplished with this filter here. If you are not familiar with Python, we recommend first taking a look at one of the introductory guides such as the official Python Tutorial <sup>[1]</sup>. Also, if you are going to do any programming beyond the very basic stuff, we recommend reading up on the VTK API. The VTK website <sup>[2]</sup> has links to VTK books and online documentation. For reference, you may need to look at the VTK class documentation <sup>[3]</sup>. There is also more information about the Programmable Filter and some good recipes on the ParaView Wiki (Python\_Programmable\_Filter).

## Basic Use

Requirements:

1. You are applying Programmable Filter to a "simple" dataset and not a composite dataset such as multi-block or AMR.
2. You have NumPy installed.

The most basic reason to use the Programmable Filter is to add a new array by possibly deriving it from arrays in the input. This can be achieved by using the Python Calculator. One reason to use the Programmable Filter instead may be that the calculation is more involved and trying to do it in one expression may be difficult. Another reason may be that you need access to a program flow construct such as if or for. In any case, the Programmable Filter can be used to do everything the Calculator does and more.

**Note:** Since what we describe here builds on some of the concepts introduced in the Python Calculator section, please read it first if you are not familiar with the Calculator.

If you leave the "Output Dataset Type" parameter in the default setting of "Same as Input", the Programmable Filter will copy the topology and geometry of the input to the output before calling your Python script. Therefore, if you Apply the filter without filling the script, you should see a copy of the input without any of its arrays in the output. If you also check the "Copy Arrays" option, the output will have all of the input arrays. This behavior allows you to focus on creating new arrays without worrying about the mesh.

Let's try an example. Create a Sphere source and then apply the Programmable Filter. Use the following script.

```
normals = inputs[0].PointData['Normals']
output.PointData.append(normals[:,0], "Normals_x")
```

This should create a sphere with one array called "Normals\_x". There are a few things to note here:

- You cannot refer to arrays directly by name as in the Python Calculator. You need to access arrays using the .PointData and .CellData qualifiers.
- Unlike the Python Calculator, you have to explicitly add an array to the output using the append function. Note that this function takes the name of the array as the second argument.

You can use any of the functions available in the Calculator in the Programmable Filter. For example, the following code creates two new arrays and adds them to the output.

```
normals = inputs[0].PointData['Normals']
output.PointData.append(sin(normals[:,0]), "sin of Normals_x")
output.PointData.append(normals[:,1] + 1, "Normals_y + 1")
```



## Intermediate Use

### Mixing VTK and NumPy APIs

The examples above demonstrate how the Programmable Filter can be used as an advanced Python Calculator. However, the full power of the Programmable Filter can only be harnessed by using the VTK API. Let's start with a simple example. Create a Sphere source and apply the Programmable Filter with the following script.

```
input = inputs[0]

newPoints = vtk.vtkPoints()
numPoints = input.GetNumberOfPoints()
for i in range(numPoints):
    x, y, z = input.GetPoint(i)
    newPoints.InsertPoint(i, x, y, 1 + z*0.3)

output.SetPoints(newPoints)
```

This requires some explanation. We start with creating a new instance of `vtkPoints`.

```
newPoints = vtk.vtkPoints()
```

`vtkPoints` is a data structure that VTK uses to store the coordinates of points. Next, we loop over all points of the input and insert a new point in the output with coordinates  $(x, y, 1+z*0.3)$

```
for i in range(numPoints):
    x, y, z = input.GetPoint(i)
    newPoints.InsertPoint(i, x, y, 1 + z*0.3)
```

Finally, we replace the output points with the new points we created using the following.

```
output.SetPoints(newPoints)
```

**Note:** Python is an interpreted language and Python scripts do not execute as efficiently as compiled C++ code. Therefore, using a for loop that iterates over all points or cells may be a significant bottleneck when processing large datasets.

The NumPy and VTK APIs can be mixed to achieve good performance. Even though this may seem a bit complicated at first, it can be used with great effect. For example, the example above can be rewritten as follows.

```
from paraview.vtk.dataset_adapter import numpyTovtkDataArray

input = inputs[0]

newPoints = vtk.vtkPoints()

zs = 1 + input.Points[:,2]*0.3
coords = hstack([input.Points[:,0:2], zs])

newPoints.SetData(numpyTovtkDataArray(coords))

output.SetPoints(newPoints)
```

Even though this produces exactly the same result, it is much more efficient because we eliminated the for loop. Well, we didn't really eliminate it. We rather moved it from Python to C. Under the hood, NumPy uses C and Fortran for tight loops.

If you read the Python Calculator documentation, this example is pretty straightforward except the use of `numpyVtkDataArray()`. First of all, note that we are mixing two APIs here: the VTK API and NumPy. VTK and NumPy uses different types of objects to represents arrays. The basic examples we previously used carefully hide this from you. However, once you start manipulating VTK objects using NumPy, you have to start converting objects between two APIs. Note that for the most part this conversion happens without "deep copying" arrays, e.g. copying the raw contents from one memory location to another. Rather, pointers are passed between VTK and NumPy whenever possible.

The `dataset_adapter` provides two methods to do the conversions described above:

- `vtkDataArrayToVTKArray`: This function create a NumPy compatible array from a `vtkDataArray`. Note that `VTKArray` is actually a subclass of `numpy.matrix` and can be used anywhere `matrix` can be used. This function always copies the pointer and not the contents. **Important:** You should not directly change the values of the resulting array if the argument is an array from the input.
- `numpyVtkDataArray`: Converts a NumPy array (or a `VTKArray`) to a `vtkDataArray`. This function copies the pointer if the argument is a contiguous array. There are various ways of creating discontinuous arrays with NumPy including using `hstack` and `striding`. See NumPy documentation for details.

## Multiple Inputs

Like the Python Calculator, the Programmable Filter can accept multiple inputs. First select two or more pipeline objects in the pipeline browser and then apply the Programmable Filter. Then each input can be accessed using the `inputs[]` variable. Note that if the Output Dataset Type is set to Same as Input, the filter will copy the mesh from the first input to the output. If Copy Arrays is on, it will also copy arrays from the first input. As an example, the following script compares the Pressure attribute from two inputs using the difference operator.

```
output.append(inputs[1].PointData['Pressure'] -
inputs[0].PointData['Pressure'], "difference")
```

## Dealing with Composite Datasets

So far, none of the example we used apply to multi-block or AMR datasets. When talking about the Python Calculator, we did not have to differentiate between simple and composite datasets. This is because the calculator loops over all of the leaf blocks of composite datasets and applies the expression to each one. Therefore, inputs in an expression are guaranteed to be simple datasets. On the other hand, the Programmable Filter does not perform this iteration and passes the input, composite or simple, as it is to the script. Even though this makes basic scripting harder for composite datasets, it provides enormous flexibility.

The only thing that you have to know to work with composite datasets is how to iterate over them to access the leaf nodes. Let's start with a simple example.

```
for block in inputs[0]:
    print block
```

Here we iterate over all of the non-NULL leaf nodes (i.e. simple datasets) of the input and print them to the Output Messages console. Note that this will work only if the input is multi-block or AMR.

Note that when Output Dataset Type is set to "Same as Input", the Programmable Filter will copy composite dataset to the output - it will copy only the mesh unless Copy Arrays is on. Therefore, you can also iterate over the output. A simple trick is to turn on Copy Arrays and then use the arrays from the output when generating new ones. Below is

an example. We use the can.ex2 file from the ParaView testing dataset collection.

```
def process_block(block):
    displ = block.PointData['DISPL']
    block.PointData.append(displ[:,0], "displ_x")

for block in output:
    process_block(block)
```

Alternatively, you can use the MultiCompositeDataIterator to iterate over the input and output block simultaneously. The following is equivalent to the previous example.

```
def process_block(input_block, output_block):
    displ = input_block.PointData['DISPL']
    output_block.PointData.append(displ[:,0], "displ_x")

from paraview.vtk.dataset_adapter import MultiCompositeDataIterator
iter = MultiCompositeDataIterator([inputs[0], output])

for input_block, output_block in iter:
    process_block(input_block, output_block)
```

## Advanced (but read it anyway)

### Changing Output Type

So far, all of the examples we discussed depended on that the output type being the same as input and that the Programmable Filter copied the input mesh to the output. If we set the output type to something other than Same as Input, we are on our own: the Programmable Filter will create an empty output of the type we specified but will not copy any information. Even though it may be more work, this gives us a lot of flexibility. Since we are approaching the realm of full-blown VTK filter authoring, we will pick a very simple example here. If you are already familiar with VTK API, you will realize that this is a great way of prototyping VTK filters. If you are not, reading up on VTK is a good idea.

Create a Wavelet source, apply a Programmable Filter, set the output type to vtkTable and use the following script:

```
rtdata = inputs[0].PointData['RTData']

output.RowData.append(min(rtdata), 'min')
output.RowData.append(max(rtdata), 'max')
```

Here, we added two columns to the output table. First one has one value - minimum of RTData - and the second one has the maximum of RTData. When you apply this filter, the output should automatically be shown in a Spreadsheet view. We could also use this sort of script to chart part of our input data. For example, the output of the following script can be display as a line chart.

```
rtdata = inputs[0].PointData['RTData']
output.RowData.append(rtdata, 'rtdata')
```

Changing the output type is also often necessary when using VTK filters within the script. We will demonstrate how to do that in the following section.

## Dealing with Structured Data Output

Let's say that we have a curvilinear grid. A good example is bluntfin.vts from the ParaView testing data. We would like to volume render a subset of this grid. As of 3.10, ParaView does not support volume rendering of curvilinear grids so we have two choices:

- Resample to an image data
- Convert to unstructured grid

Here we will demonstrate how to resample to an image data using the Programmable Filter. Of course, this can be accomplished using the Resample with Dataset filter but it is a good example nevertheless. Start with loading bluntfin.vts, then apply the Programmable Filter. Make sure to set the output type to vtkImageData. Here is the script:

```
pinput = vtk.vtkImageData()
pinput.SetExtent(0, 10, 0, 10, 0, 10)
pinput.SetOrigin(0, 1, 0)
pinput.SetSpacing(0.5, 0.5, 0.5)

probe = vtk.vtkProbeFilter()
probe.SetInput(pinput)

input_copy = inputs[0].NewInstance()
input_copy.UnRegister(None)
input_copy.ShallowCopy(inputs[0].VTKObject)

probe.SetSource(input_copy)
probe.Update()

output.ShallowCopy(probe.GetOutput())
```

**Note:** See the next section for details about using a VTK filter within the Programmable Filter.

If you already applied, you may notice that the output looks much bigger than it should be. This is because we are missing an important piece. We need to use the following as the RequestInformation script:

```
from paraview.util import SetOutputWholeExtent

SetOutputWholeExtent(self, [0, 10, 0, 10, 0, 10])
```

This requires a bit of explanation. VTK expects that all data sources and filters that produce structured data (rectilinear or curvilinear grids) to provide meta data about the logical extents of the output dataset before full execution. Thus the RequestInformation is called by the Programmable Filter before execution and is where you should provide this meta-data. This is not required if the filter is simply copying the mesh as the meta-data would have been provided by another pipeline object upstream. However, if you are using the Programmable Filter to produce a structured data with a different mesh than the input, you need to provide this information.

You are probably wondering what the RequestUpdateExtent script does. This script can be used to augment the request that propagates upstream before execution. This is used to ask for a specific data extent, for example. This is an advanced concept and is not discussed further here.

## Using VTK Filters with Programmable Filter

The previous example demonstrated how you can use a VTK filter (vtkProbeFilter in this case) from with the Programmable Filter. We will explain that example in more detail here. Here it is again.

```
pinput = vtk.vtkImageData()
pinput.SetExtent(0, 10, 0, 10, 0, 10)
pinput.SetOrigin(0, 1, 0)
pinput.SetSpacing(0.5, 0.5, 0.5)

probe = vtk.vtkProbeFilter()
probe.SetInput(pinput)

input_copy = inputs[0].NewInstance()
input_copy.UnRegister(None)
input_copy.ShallowCopy(inputs[0].VTKObject)

probe.SetSource(input_copy)
probe.Update()

output.ShallowCopy(probe.GetOutput())
```

There are two important tricks to use a VTK filter from another VTK filter. First, do not directly set the input to the outer filter as the input of the inner filter. It is difficult to explain why without getting into VTK pipeline mechanics so please take our word for it. Instead, make a shallow copy as follow.

```
input_copy = inputs[0].NewInstance()
input_copy.UnRegister(None)
input_copy.ShallowCopy(inputs[0].VTKObject)
```

The UnRegister() call is essential to avoid memory leaks.

The second trick is to use ShallowCopy() to copy the output of the internal filter to the output of the outer filter as follows.

```
output.ShallowCopy(probe.GetOutput())
```

This should be enough to get you started. There are a large number of VTK filters so it is not possible to describe them here. Refer to the VTK documentation for more information.

## References

- [1] <http://docs.python.org/tutorial/>
- [2] <http://www.vtk.org>
- [3] <http://www.vtk.org/doc/release/5.6/html/>

# Calculator

## Basics

The Calculator Filter can be used to calculate derived quantities from existing attributes. The main parameter of the Calculator is an expression that describes how to calculate the derived quantity. You can enter this expression as free-form text or using some of the shortcuts (buttons and menus provided). There are some "hidden" expressions for which there are no buttons. Operands that are accessible only by typing in the function name include:

- **min(expr1, expr2)**  
returns the lesser of the two scalar expressions
- **max(expr1, expr2)**  
returns the greater of the two scalar expressions
- **cross(expr1, expr2)**  
returns the vector cross product of the two vector expressions
- **sign(expr)**  
returns -1, 0 or 1 depending if the scalar expression is less than, equal to or greater than zero respectively
- **if(condition,true\_expression,false\_expression)**  
evaluates the conditional expression and then evaluates and returns one of the two expressions
- **>** : numerical "greater than" conditional test
- **<** : numerical "less than" conditional test
- **=** : numerical "equal to" conditional test
- **&** : boolean "and" test conjunction
- **|** : boolean "or" test conjunction

Here we describe various options available within the Calculator.

**Note:** We recommend that you use the Python Calculator instead of Calculator if possible. The Python Calculator is more flexible, has more functions and is more efficient. However, it requires that ParaView is compiled with Python support and that NumPy is installed.

Let's start with a simple example. Create a Wavelet source and then apply the Calculator using "1" as the expression.

**Note:** You can enter an expression by clicking in the expression entry box and typing. This should create an point array called "Result" in the output. A few things to note:

- The Calculator copies the input mesh to the output. It is possible to have the calculator to change the point coordinates. We will discuss this later.
- The expression is calculated for each element in the output point or cell data (depending on the Attribute Mode).

Next, change the expression to be "5 \* RTData" and the Result Array Name to be "5 times rtdata" (without the quotes). If you change to surface representation and color by the new array, you will notice that the filter calculated "5 \* RTData" at each point.

The main use case for the Calculator is to utilize one or more input arrays to calculate derived quantities. The Calculator can either work on point centered attributes or cell centered attributes (but not both). In order to help enter the names of the input arrays, the Calculator provides two menus accessible through the "Scalars" and "Vectors" buttons. If you select an array name from either menus, it will be inserted to the expression entry box at the cursor location. You can also use the other buttons to enter any of the functions available to the Calculator. Most of these functions are obvious. We will cover the ones that are not below.

## Working with Vectors

Let's start with an example. Create a Wavelet source then apply the Random Vectors filter. Next, apply the Calculator. Now look at the Scalars and Vectors menus on the Object Inspector panel. You will notice that BrownianVectors shows up under Vectors, whereas BrownianVectors\_X, \_Y and \_Z show up under scalars. The Calculator allows access to individual components of vectors using this naming convention. So if we use BrownianVectors\_X as the expression, the Calculator will extract the first component of the BrownianVectors attribute. All of the Calculator's functions are applicable to vectors. Most of these functions treat the vector attributes same as scalars, mainly apply the same functions to all components of all elements. However, the following functions work only on vectors:

- **v1 . v2** : Dot product of two vectors. Returns a scalar.
- **norm** : Creates a new array which contains normalized versions of the input vectors.
- **mag** : Returns the magnitude of input vectors.

You may have noticed that 4 calculator buttons on the Object Inspector are not actually functions. Clear is straightforward. It cleans the expression entry box. iHat, jHat and kHat on the other hand are not as clear. These represent unit vectors in X, Y and Z directions. They can be used to construct vectors from scalars. Let's consider a simple but common use case. We want to set the Z component of BrownianVectors from the previous example to 0. The expression to do that is "BrownianVectors\_X \*iHat+BrownianVectors\_Y\*jHat+0\*kHat". This expression multiplies the X unit vector with the X component of the input vector, the Y unit vector with the Y component, and the Z unit vector with 0 and add them together. You can use this sort of expression to create vectors from individual components of a vector if the reader loaded them separately, for example. **Note:** We didn't really need the 0\*kHat bit. It was for demonstration.

## Working with Point Coordinates

You may have noticed that one point-centered vector and its 3 components are always available in the Calculator. This vector is called "coords" and represents the point coordinates. You can use this array in your expression like any other array. For example, in our example above, we could use "mag(coords)\*RTData" to scale RTData with the distance of the point to the origin.

It is also possible to change the coordinates of the mesh by checking the "Coordinate Results" box. Note that this does not work for rectilinear grids (uniform and non-uniform) since their point coordinates cannot be adjusted one-by-one. Since our previous examples used a uniform rectilinear grid, we cannot use them. Instead, start with the Sphere source, then use this expression: "coords+2\*iHat". Make sure to check the "Coordinate Results" box. The output of the Calculator should be a shifted version of the input sphere.

## Dealing with Invalid Results

Certain functions are not applicable to certain arguments. For example, sqrt() works only on positive numbers since the calculator does not support complex numbers. Unless the "Replace invalid results" option is turned on, an expression that tries to evaluate the square root of a negative number will return an error such as this:

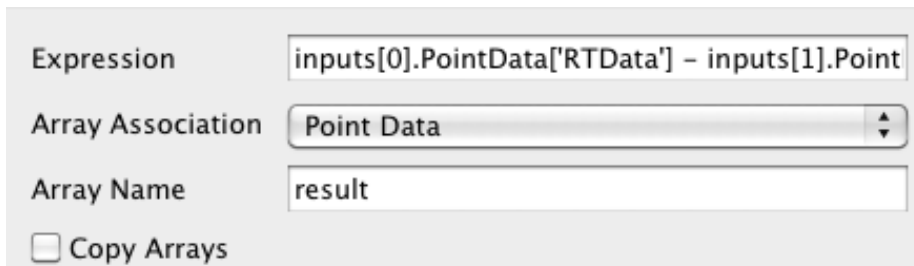
```
ERROR: In /Users/berk/Work/ParaView/git/VTK/Common/vtkFunctionParser.cxx, line 697  
vtkFunctionParser (0x128d97730): Trying to take a square root of a negative value
```

However, if you turn on the "Replace invalid results" option, the calculator will silently replace the result of the invalid expression with the value specified in "Replacement value". Note that this will happen only when the expression result in an invalid result so some of the output points (or cells) may have the Replacement Value whereas others may have valid results.

# Python Calculator

---

## Introduction



The image shows the Python Calculator widget interface. It has four main sections: 'Expression' with a text input field containing 'inputs[0].PointData['RTData'] - inputs[1].Point'; 'Array Association' with a dropdown menu set to 'Point Data'; 'Array Name' with a text input field containing 'result'; and a 'Copy Arrays' checkbox which is currently unchecked.

The Python Calculator is a ParaView filter that processes one or more input arrays based on an expression provided by the user to produce a new output array. The parameters of the filter include the expression, the association of the output array (Point or Cell Data), the name of output array and a toggle that controls whether the input arrays are copied to the output. In this document, we introduce the use of the Python Calculator and provide a list of functions available to the user.

Note that the Python Calculator depends on Python and NumPy. All ParaView binaries distributed by Kitware are built with these to enable the calculator. If you have built ParaView yourself, you have to make sure that NumPy is installed and that PARAVIEW\_ENABLE\_PYTHON is turned on when configuring the ParaView build.



## Basic Tutorial

Start by creating a Sphere source and applying the Python Calculator to it. As the first expression, use the following and apply:

```
5
```

This should create an array name "result" in the output point data. Note that this is an array that has a value of 5 for each point. When the expression results in a single value, the calculator will automatically make an constant array. Next, try the following:

```
Normals
```

Now the "result" array should be the same as the input array Normals. As described in detail later, various functions are available through the calculator. For example, the following is a valid expression.

```
sin(Normals) + 5
```

It is very important to note that the Python Calculator has to produce one value per point or cell depending on the Array Association parameter. Most of the functions described here apply individually to all point or cell values and produce an array as the same dimensions as the input. However, some of them (such as min() and max()) produce single values.

## Accessing Data

There are several ways of accessing input arrays within expressions. The simplest way is to access it by name:

```
sin(Normals) + 5
```

This is equivalent to:

```
sin(inputs[0].PointData['Normals']) + 5
```

The example above requires some explanation. Here inputs[0] refer to the first input (dataset) to the filter. Python Calculator can accept multiple inputs. Each input can be accessed as inputs[0], inputs[1], ... You can access the point or cell data of an input using the .PointData or .CellData qualifiers. You can then access individual arrays within the point or cell data containers using the [] operator. Make sure to use quotes or double-quotes around the array name. NOTE: Arrays that have names with certain characters (such as space, +, -, \*, /) in their name can only be accessed using this method.

Certain functions apply directly on the input mesh. These filters expect an input dataset as argument. For example,

```
area(inputs[0])
```

For data types that explicitly define the point coordinates, you can access the coordinates array using the .Points qualifier. The following extracts the first component of the coordinates array:

```
inputs[0].Points[:,0]
```

Note that certain data types, mainly image data (uniform rectilinear grid) and rectilinear grid, point coordinates are defined implicitly and cannot be accessed as an array.

## Comparing Multiple Datasets

As described above, the Python Calculator can accept multiple inputs. Using this functionality, you can compare multiple datasets. There are a few things to note:

- Python Calculator will always copy the mesh from the first input to its output,
- All operations are applied point by point. In most cases, this requires that the input meshes (topology and geometry) are the same. At the least, it requires the the inputs have the same number of points or cells
- In parallel execution mode, the inputs have to be distributed exactly the same way across processes

For example, to compare the temperature of 2 datasets, you can select both inputs, apply the Python Calculator and then use the following expression.

```
inputs[0].PointData['temperature'] - inputs[1].PointData['temperature']
```

## Basic Operations

The Python calculator supports all of the basic arithmetic operations using the +, -, \* and / operators. These are always applied element-by-element to point and cell data including scalars, vectors and tensors. These operations also work with single values. For example, the following adds 5 to all components of all Normals.

```
Normals + 5
```

The following adds 1 to the first component, 2 to the second component and 3 to the third component:

```
Normals + [1, 2, 3]
```

This is specially useful when mixing functions that return single values. For example, the following normalizes the Normals array:

```
(Normals - min(Normals)) / (max(Normals) - min(Normals))
```

A common use case in a calculator is to work on one component of an array. This can be accomplished with the following:

```
Normals[:, 0]
```

The expression above extracts the first component of the Normals vector. Here, : is a placeholder for "all elements". One element can be extracted by replacing : with an index. For example, the following creates an constant array from the first component of the normal of the first point:

```
Normals[0, 0]
```

Whereas the following, assigns the normal of the first point to all points:

```
Normals[0, :]
```

It is also possible to merge multiple scalars into on array. For this, we can use the `hstack()` function:

```
hstack([velocity_x, velocity_y, velocity_z])
```

Note the use of square brackets ([]).

Under the cover, the Python Calculator uses NumPy. All arrays in the expression are compatible with NumPy arrays and can be used where NumPy arrays can be used. For more information on what you can do with these arrays, consult with the NumPy book, which can be downloaded here <sup>[1]</sup>.

## Functions

The following is a list of functions available in the Python Calculator. Note that this list is partial since most of the NumPy and SciPy functions can be used in the Python Calculator. Many of these functions can take single values or arrays as argument.

**abs (x)** : Returns the absolute value(s) of x.

**add (x, y)**: Returns the sum of two values. x and y can be single values or arrays. Same as x+y.

**area (dataset)** : Returns the surface area of each cell in a mesh.

**aspect (dataset)** : Returns the aspect ratio of each cell in a mesh.

**aspect\_gamma (dataset)** : Returns the aspect ratio gamma of each cell in a mesh.

**condition (dataset)** : Returns the condition number of each cell in a mesh.

**cross (x, y)** : Return the cross product for two 3D vectors from two arrays of 3D vectors.

**curl (array)**: Returns the curl of an array of 3D vectors.

**divergence (array)**: Returns the divergence of an array of 3D vectors.

**divide (x, y)**: Element by element division. x and y can be single values or arrays. Same as x/y.

**det (array)** : Returns the determinant of an array of 2D square matrices.

**determinant (array)** : Returns the determinant of an array of 2D square matrices.

**diagonal (dataset)** : Returns the diagonal length of each cell in a dataset.

**dot (a1, a2)**: Returns the dot product of two scalars/vectors of two array of scalars/vectors.

**eigenvalue (array)** : Returns the eigenvalue of an array of 2D square matrices.

**eigenvector (array)** : Returns the eigenvector of an array of 2D square matrices.

**exp (x)**: Returns power(e, x).

**global\_max(array)**: Returns the maximum value of an array of scalars/vectors/tensors among all process. Not yet supported for multi-block and AMR datasets.

**global\_mean (array)** : Returns the mean value of an array of scalars/vectors/tensors among all process. Not yet supported for multi-block and AMR datasets.

**global\_min(array)**: Returns the minimum value of an array of scalars/vectors/tensors among all process. Not yet supported for multi-block and AMR datasets.

**gradient(array)**: Returns the gradient of an array of scalars/vectors.

**inv (array)** : Returns the inverse an array of 2D square matrices.

**inverse (array)** : Returns the inverse of an array of 2D square matrices.

**jacobian (dataset)** : Returns the jacobian of an array of 2D square matrices.

**laplacian (array)** : Returns the jacobian of an array of scalars.

**ln (array)** : Returns the natural logarithm of an array of scalars/vectors/tensors.

**log (array)** : Returns the natural logarithm of an array of scalars/vectors/tensors.

**log10 (array)** : Returns the base 10 logarithm of an array of scalars/vectors/tensors.

**max (array)**: Returns the maximum value of the array as a single value. Note that this function returns the maximum within a block for AMR and multi-block datasets, not across blocks/grids. Also, this returns the maximum within each process when running in parallel.

**max\_angle (dataset)** : Returns the maximum angle of each cell in a dataset.

**mag (a)** : Returns the magnigude of an array of scalars/vectors.

**mean (array) :** Returns the mean value of an array of scalars/vectors/tensors.

**min (array) :** Returns the minimum value of the array as a single value. Note that this function returns the minimum within a block for AMR and multi-block datasets, not across blocks/grids. Also, this returns the minimum within each process when running in parallel.

**min\_angle (dataset) :** Returns the minimum angle of each cell in a dataset.

**mod (x, y):** Same as remainder (x, y).

**multiply (x, y):** Returns the product of x and y. x and y can be single values or arrays. Note that this is an element by element operation when x and y are both arrays. Same as  $x * y$ .

**negative (x):** Same as  $-x$ .

**norm (a) :** Returns the normalized values of an array of scalars/vectors.

**power (x, a):** Exponentiation of x with a. Here both x and a can either be a single value or an array. If x and y are both arrays, a one-by-one mapping is used between two arrays.

**reciprocal (x):** Returns  $1/x$ .

**remainder (x, y):** Returns  $x - y*\text{floor}(x/y)$ . x and y can be single values or arrays.

**rint (x):** Rounds x to the nearest integer(s).

**shear (dataset) :** Returns the shear of each cell in a dataset.

**skew (dataset) :** Returns the skew of each cell in a dataset.

**square (x):** Returns  $x*x$ .

**sqrt (x):** Return square root of x.

**strain (array) :** Returns the strain of an array of 3D vectors.

**subtract (x, y):** Returns the difference between two values. x and y can be single values or arrays. Same as  $x - y$ .

**surface\_normal (dataset) :** Returns the surface normal of each cell in a dataset.

**trace (array) :** Returns the trace of an array of 2D square matrices.

**volume (dataset) :** Returns the volume normal of each cell in a dataset.

**vorticity(array):** Returns the vorticity/curl of an array of 3D vectors.

**vertex\_normal (dataset) :** Returns the vertex normal of each point in a dataset.

## Trigonometric Functions

Below is a list of supported trigonometric functions.

**sin (x)**

**cos (x)**

**tan (x)**

**arcsin (x)**

**arccos (x)**

**arctan (x)**

**hypot (x1, x2)**

**sinh(x)**

**cosh (x)**

**tanh (x)**

**arcsinh (x)**

arccosh (x)

arctanh (x)

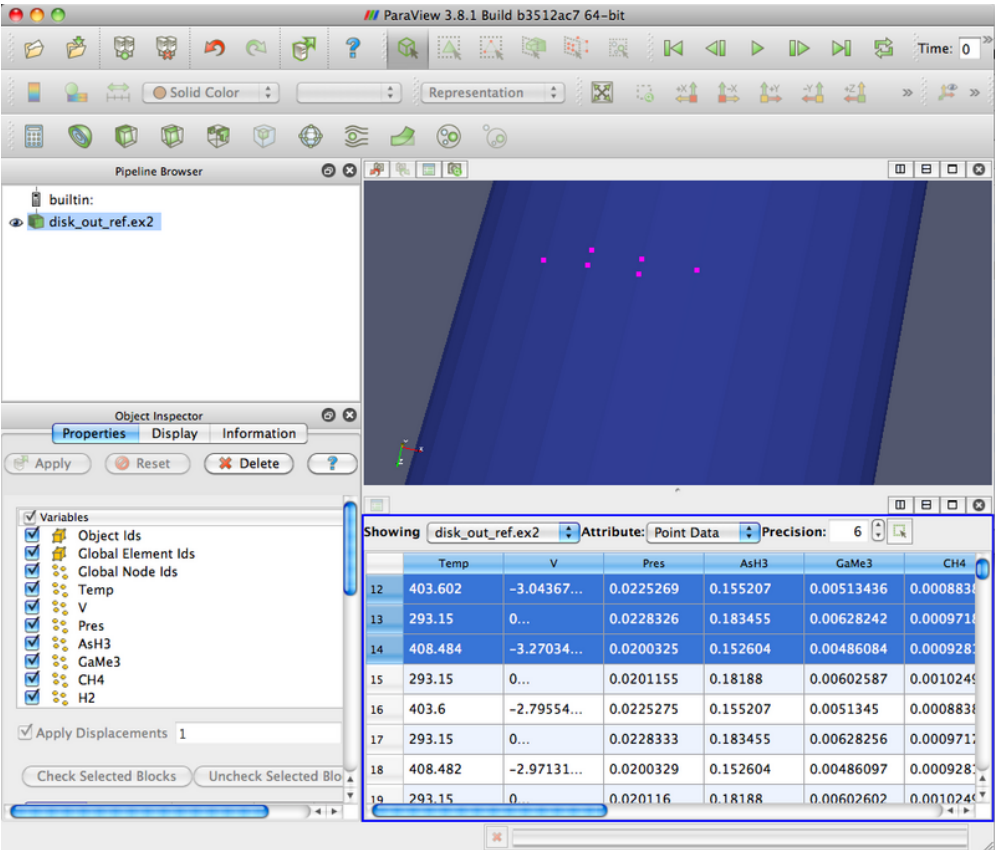
References

[1] <http://www.tramy.us/guidetoscipy.html>

Spreadsheet View

In several cases it’s very useful to look at the raw dataset. This is where the Spreadsheet view comes in handy. Spreadsheet View allows users to explore the raw data in a spreadsheet-like display. Users can inspect cells and points and data values associated with these using this view. This makes it a very useful for drilling down into the data.

Spreadsheet View, as the name suggests is a view i.e. users can create one by splitting the main-view frame in the application client-area. Refer to the chapter on Views for details on views. Spreadsheet View can only show one dataset at a time. However, users can use ParaView’s multi-view capabilities to create multiple spreadsheet views for inspecting different datasets.



## Inspecting Large Datasets

Spreadsheet View is a client-only view i.e. it delivers the necessary data to the client, when running in client-server mode. Additionally, it is not available through Python scripting (except when running through the Python shell provided by the ParaView application) or batch scripting. Furthermore, when running on a tile-display, the area covered by the spreadsheet on the client simply shows up as a blank region on the tiles.

Unlike complex views like the 3D render view, that can generate the renderings on the server and simply deliver images to the client, the spreadsheet view requires that the data is available on the client. This can be impractical when inspecting really large datasets since the client may not even have enough memory, even if we assume infinite bandwidth. To address such issues, the Spreadsheet View streams the data to the client, only fetching the data for the row currently visible in the viewport.

## Double Precision

Using the precision spin-box in the view header, users can control the precision for floating point numbers. The value determines the number of significant digits after the decimal point.

## Selection with Spreadsheet View

Spreadsheet View, in many ways, behaves like typical spreadsheet applications. One can scroll, select rows using mouse clicks, arrow keys and modifiers like Shift and Ctrl keys, sort columns by clicking on the header. Additionally, one can double-click on a column header to toggle the maximization of a column for better readability.

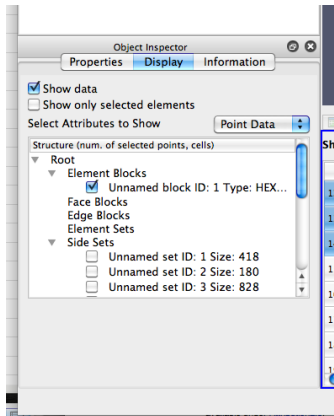
On selecting rows the corresponding cells or points will get selected and ParaView will highlight those in other views, such as the 3D view. Conversely, when one makes a selection in the 3D View or the chart views, one the rows corresponding to the selected cells/points will be highlighted in the spreadsheet view. Of course, for the view to highlight the selection, the selected dataset must be the one that's being shown in the spreadsheet view. To make it easier to inspect just the selected elements, one check the "Show only selected" button on the view header.

When in "Show only Selected" mode, one can no longer create selections on the Spreadsheet view. One has to use the other views to make the selection and the Spreadsheet view will automatically update to show the details for the items that got selected.

Spreadsheet View can show data associated with cells, points or even field data. To choose what attribute the view shows, one can use the attribute-type combo-box. The selection created by the view depends on the attribute-type, i.e. when if user selects in the view when attribute type is "Points", points in the dataset will be selected. The spreadsheet view also performs selection conversions if possible, i.e. if you select a cell in the 3D view, however you the spreadsheet view is setup to show points, then the view will highlight the points that form the selected cell.

The Spreadsheet View may add several additional data columns which may not be present in the your actual data. These data-columns are either derived information such as the (i, j, k) co-ordinates for structured data or provide additional information about the data, e.g. block index for composite datasets, or provide additional information about the data distribution such as process-id when connected to a parallel server (pvserver or pvdataserver).

## Working with Composite Datasets



Spreadsheet view works seamlessly with different kinds of dataset types including composite datasets such as multi-block datasets or AMR datasets. When dealing with composite datasets, the view shows one block at a time. Users can choose the block to inspect by using the display tab.

## Selection

---

Selection is the mechanism for identifying a subset of a dataset by using user specified criteria. This subset can be a set of point or cells or a block of composite dataset. This functionality allows users to focus on a smaller subset that is important. For example, the elements of a finite-element mesh that have pressure above a certain threshold can be identified very easily using the threshold selection. Furthermore, this selection can be converted to a set of global element ids in order to plot the attribute values of those elements over time.

In this chapter we discuss the mechanism for creating selections using **Views** and **Selection Inspector**. In the following chapter, we detail another powerful and flexible mechanism of creating selection using queries.

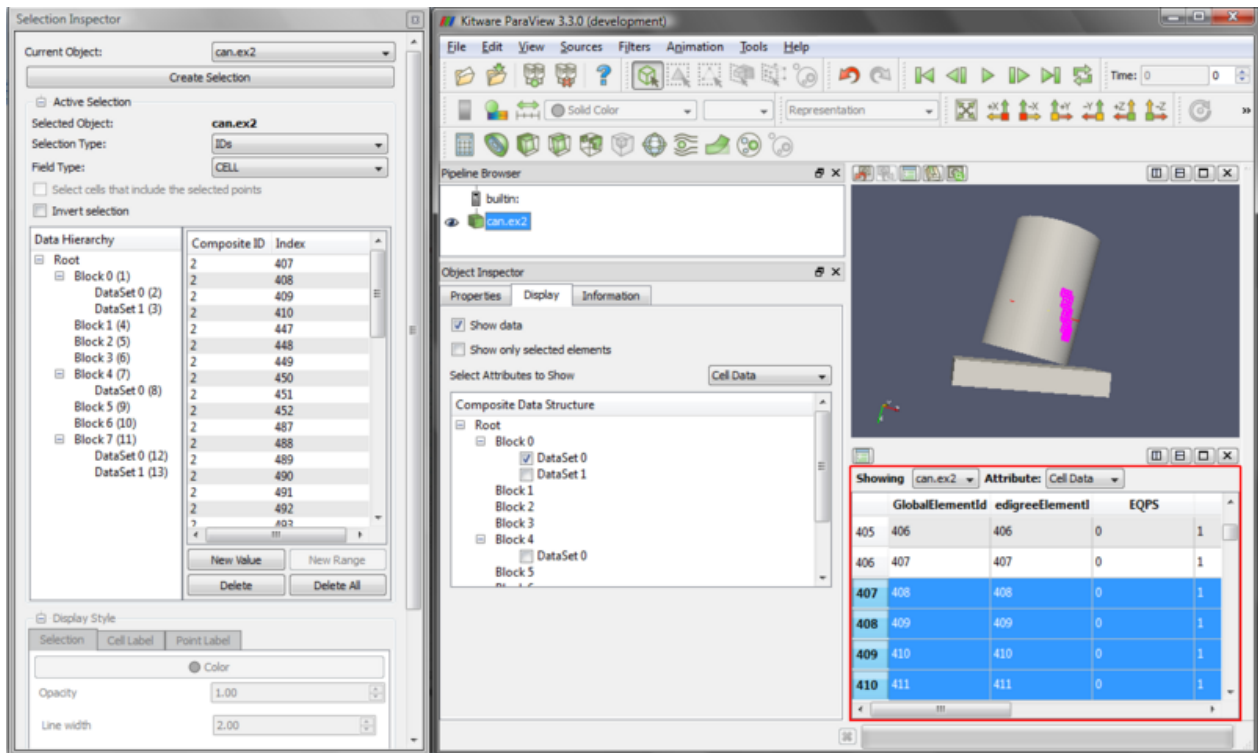
ParaView supports a single active selection. This selection is associated with a data source (here data source refers to any reader, source or filter) and is shown in every view that displays the data source's output. This article uses a use-case driven approach to demonstrate how this selection can be described and used. In the next section, we introduce the main GUI components that are used in the article. The subsequent sections address different use cases.

## Selection Inspector

ParaView provides a selection inspector (referred to simply as the inspector in this article) to inspect and edit the details about the active selection. One can toggle the inspector visibility from the View menu. The inspector can be used to create a new active selection, view/edit the properties of the active selection as well as change the way the selection is displayed in the 3D window e.g. change the color, show labels etc.

We will look at each of these options as we try to explore the different selection types in ParaView.

## Spreadsheet View



Spreadsheet View provides data exploration capabilities. One of the common complaints many users have is not being able to look at the raw data directly. Spreadsheet view provides exactly that. It allows the user to look at the raw cell data, point data or field data associated with a dataset. For more details on how to use the Spreadsheet View, please refer to Spreadsheet View Chapter.

## Create a Selection

In this section we will discuss different ways of creating a selection.

### Select cells/points on the Surface

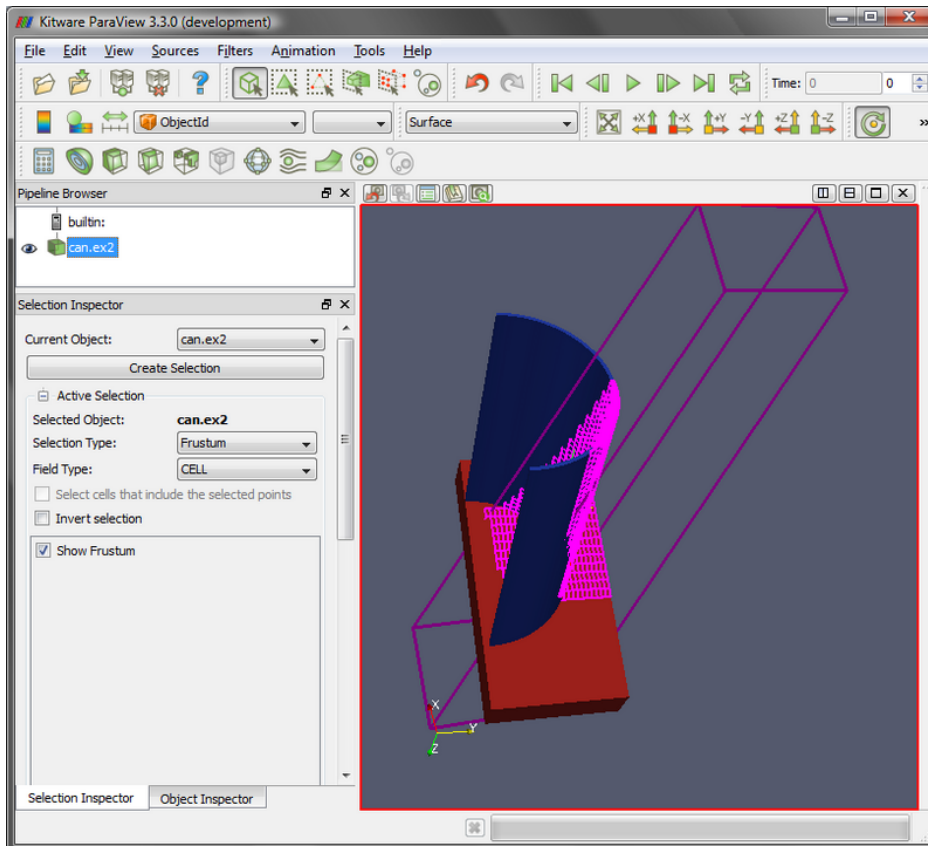
One of the simplest use-cases is to select cells or points on the surface of the dataset. It is possible to select surface cells by drawing a rubber-band on the 3D view. With the 3D view showing the dataset active, click on Select Cells (or Points) On in the Selection Controls toolbar or under the Edit menu (you can also use the 'S' key a shortcut for 'Select Cells On'). This will put ParaView into a selection mode. In this mode, click and drag over the surface of the dataset in the active view to select the cells (or points) on the surface. If anything was selected, it will be highlighted in all the views showing the data and the source producing the selected dataset will become active in the Pipeline Browser. ParaView supports selecting only one source at a time. Hence even if you draw the rubber band such that it covers data from multiple sources, only one of them will be selected (the one that has the largest number of selected cells or points).

As mentioned earlier, when data from a source is selected, all the views displaying the data show the selection. This includes spreadsheet view as well. If the spreadsheet view will show cell or point attributes of the selected data, then it will highlight the corresponding rows. When selecting points, the spreadsheet view will show the selection only if point attributes are being displayed. When selecting cells, it will highlight the cells in the cell attribute mode, and highlight the points forming the cells in the point attribute mode. For any decent sized dataset, it can be a bit tricky to locate the selected rows. In that case, the Show only selected elements on the display tab can be used to hide all the rows that were not selected.



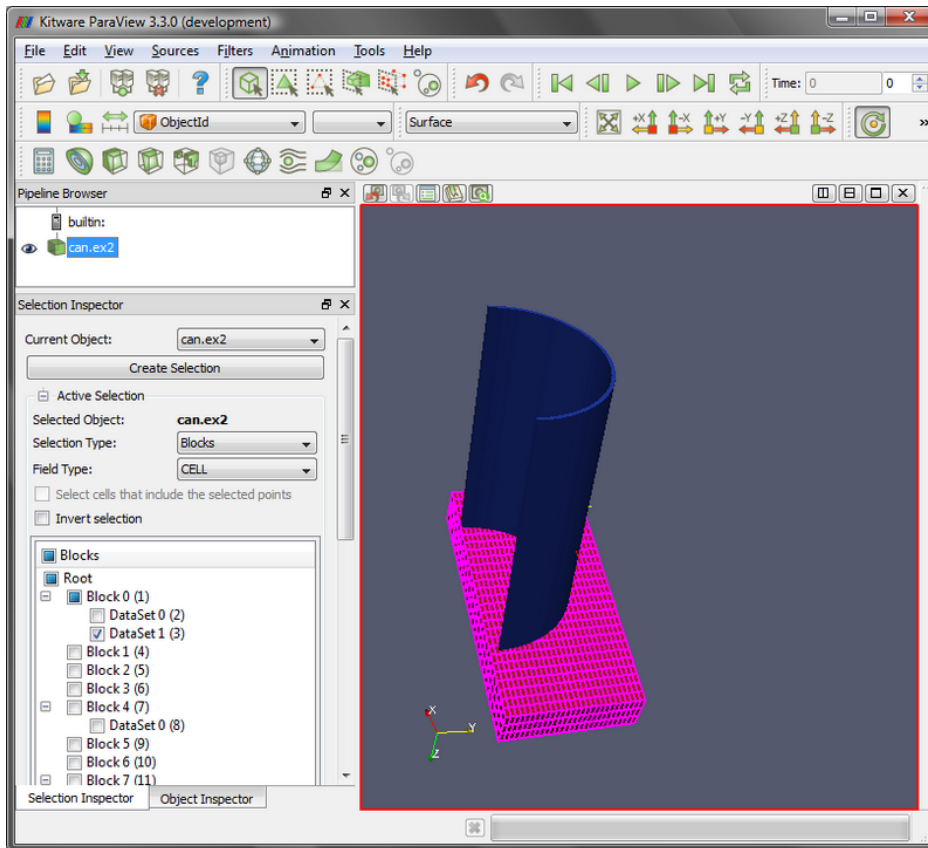
When selecting cells (or points) on the surface, ParaView determines the cell (or point) ids for each of the cell (or point) rendered within the selection box. The selection is simply the ids for cells (or points) thus determined.

## Select cells/points using a Frustum



This is similar to selecting on the surface except that instead of selecting the cells (or points) on the surface of the dataset, it selects all cells (or points) that lie within the frustum formed by extruding the rectangular rubber band drawn on the view into 3D space. To perform a frustum selection, we use Select Cells (or Points) Through in the Selection Controls toolbar or under the Edit menu. As with surface selection, the selected cells/points are shown in all the views in which the data is shown including the spreadsheet view. Unlike surface selection, the indices of the cells or points are not computed after a frustum selection. Instead ParaView performs intersections to identify the cells (or points) that lie within the frustum. Note that this selection can produce a very large selection. This may be time consuming and can increase the memory usage significantly.

## Select Blocks in a Composite Dataset



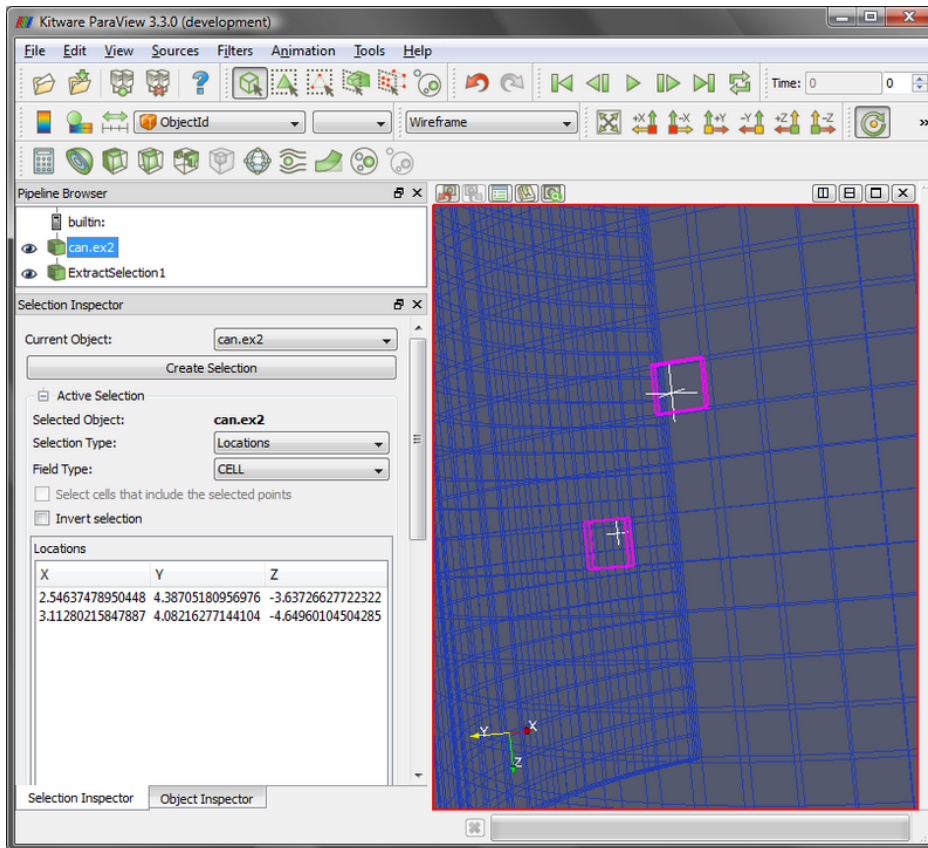
Composite datasets are multi-block or AMR (adaptive mesh refinement) datasets. In case of multi-block datasets, each block may represent different components of a large assembly e.g. tires, chassis etc. for a car dataset. Just like selecting cells or points, it is possible to select entire blocks. To enter the block selection mode use Select Block in the Selection Controls toolbar or under the Edit menu (you can also use the 'B' key as a shortcut for Select Block). Once in block selection mode, you can simply click on the block in the 3D view to select a single block or click and drag to select multiple blocks. When a block is selected, its surface cells will be highlighted.

## Select using the Spreadsheet View

Until now we have been looking at defining the selection on the 3D view. Now we will see how to create selections using the spreadsheet view. As we discussed earlier, the spreadsheet view simply shows the raw cell (point or field) data in a table. Each row represents a unique cell (or point) from the dataset. Like with any other spreadsheet application, one can select a cell (or a point) by simply clicking on the row to select. One can expand the selection using Ctrl, Shift keys while clicking. If the spreadsheet view is currently showing point attributes, then selecting on it will create a point based selection. Similarly, if it's showing cell attributes then it will create a cell based selection. Selection cannot be created when showing field attributes which are not associated with any cell or point.

We know that all views showing a selected dataset show the selection. The spreadsheet view showing the data from a source selected in the 3D view highlights the corresponding cells/points. Conversely, when we create a selection in the spreadsheet view, the corresponding cell (or point) gets highlighted in all the 3D views showing the source.

## Select using the Selection Inspector



We have seen how to use the different views for creating different types of selections. However, sometimes you may want to tweak the selection or create a new selection with a known set of cell (or point) ids or create selections based on value of any array or location etc. This is possible using the selection inspector.

Whenever a selection is created in any of the views, it becomes the active selection. The active selection is always shown in the selection inspector. For example, if you select cells on the surface of a dataset, then as shown in Figure [1], the selection inspector will show indices for the cells selected.

The selection inspector has three sections: the topmost Current Object and Create Selection are used to choose the source whose output you want to create a selection on. The Active Selection group shows the details of the active selection, if any. The Display Style group makes it possible to change the way the selection is shown in the active 3D view.

To create a new selection, choose the source whose output needs to be selected in the Current Object combo-box and then hit Create Selection. An empty selection will be created and its properties will be shown in the active selection group. Alternatively you can use any of the methods described earlier to create a selection. It will still be shown in the selection inspector.

When you select cells (or points) on the surface or using the spreadsheet view, the selection type is set to IDs. Creating a frustum selection results in a selection with the selection type set to Frustum, while selecting a block in a composite dataset creates a Block selection. Field Type indicates whether cells or points are to be selected.

In the active selection group, Selection Type indicates the type of the active selection. One can change the type by choosing one of the available options.

As shown in Figure [1], for IDs selection, the inspector lists all the selected cell or point indices. You can edit the list of ids to add or remove values. When connected to a parallel server, cell or point ids are not unique. Hence, one has to additionally specify the process number for each cell or point id. Process number -1 implies that the cell (or point) at the given index is selected on all processes. For multi-block datasets, we also need to indicate the block to which

the cell or point belongs; while for AMR datasets, we need to specify the (AMR level, index) pair.

As shown in Figure [2], for Frustum selection, currently only the Show Frustum option is available. When this option is turned on, ParaView shows the selection frustum in the 3D view. In the future, we will implement a 3D widget to modify the frustum.

As shown in Figure [3], for Block selection, the full composite tree is shown in the inspector with the selected blocks checked. Using the selection inspector one can create a selection based on thresholds for scalars in the dataset. Choose the scalar array and then add value ranges for the selected cells (or points).

Selection inspector can be used to create location based selection. When field type is CELL, cells at the indicated 3D locations will be selected. When field type is POINT, the point closest to the location within a certain threshold is selected. If Select cells that include the selected points is checked, then all the cells that contain the selected point are selected. It is possible to specify more than one location. To aid in choosing positions, one can turn the Show location widgets option on. As a result ParaView will show cross hairs in the active 3D view which can be moved interactively, as shown in Figure [4].

Selection inspector also provides a means to create global id based selection. These are similar to index based selection however since global ids are unique across all processes and blocks, one does not need to specify any additional as needed by the id based selection.

## Convert Selections

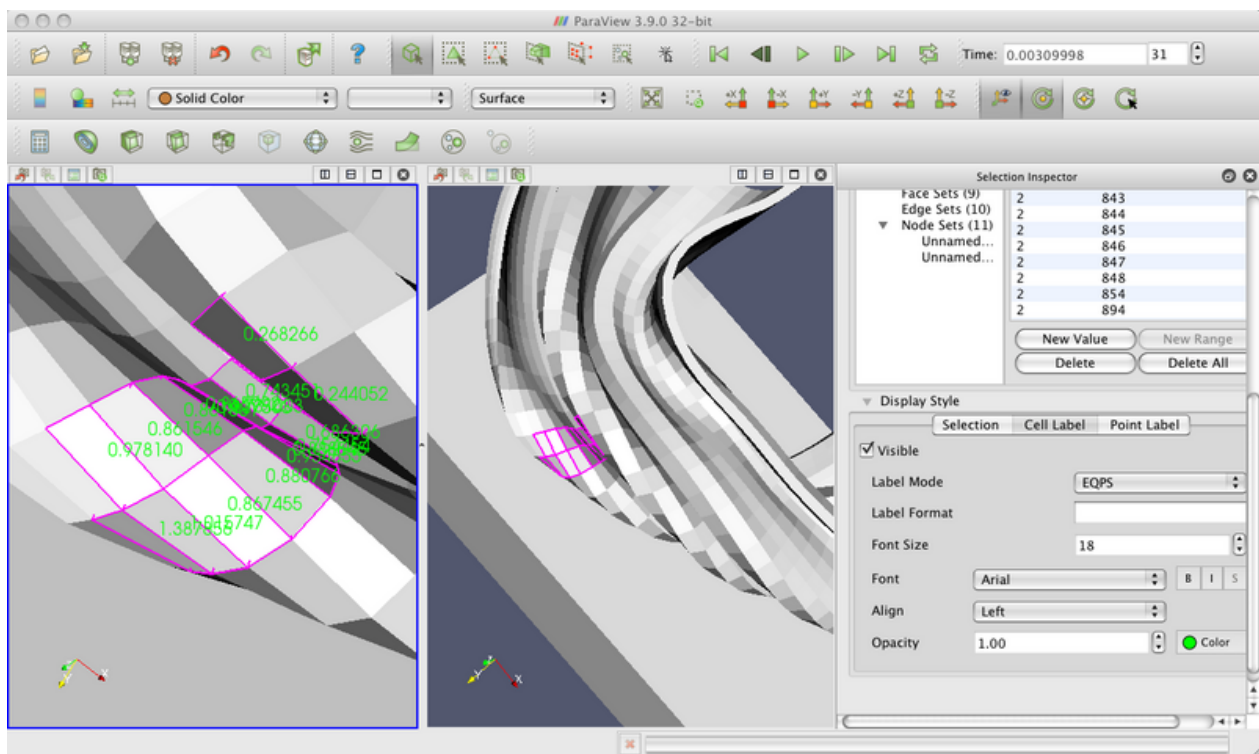
Selection inspector can also be used to convert a selection of one type to another. With some valid active selection present, if one changes the selection type then ParaView will try to convert the current selection to the new type, still preserving the cells (or points) that were selected, if possible. For example, if one create a frustum based selection and then changes the selection type to IDs, ParaView will determine the indices for all the cells (or points) that lie with the frustum and initialize the new index based selection with those indices. Note that the number of cells or points that get selected in frustum selection mode can potentially be very large; hence this conversion can be slow and memory expensive. Similarly, if the dataset provides global ids, then it is possible to convert between IDs selection and global id based selection.

It is not possible to convert between all types of selections due to obvious reasons. Conversions between id based and global id based selections, conversions from frustum to id based selections, and conversions from frustum to global id based selections are supported by the selection inspector.

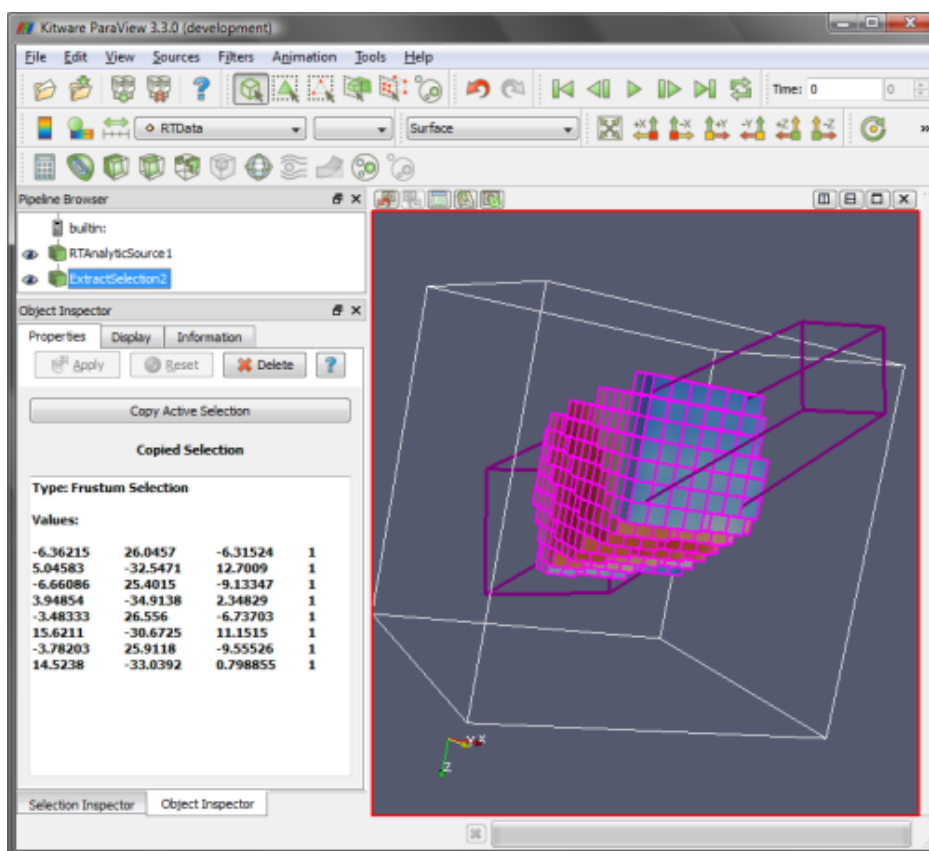
## Label Selected Cell/Points

Once an active selection is created, we can label the cells or points or their associated values in a 3D view. This can be done using the selection inspector. At the bottom of the selection inspector panel, there are two tabs Cell Label and Point Label which can be used to change the cell or point label visibility and other label attributes such color, font etc. These tabs are enabled only if the active view is a 3D view. Any changes done in the Display Style group (including the labels) only affect the active 3D view.

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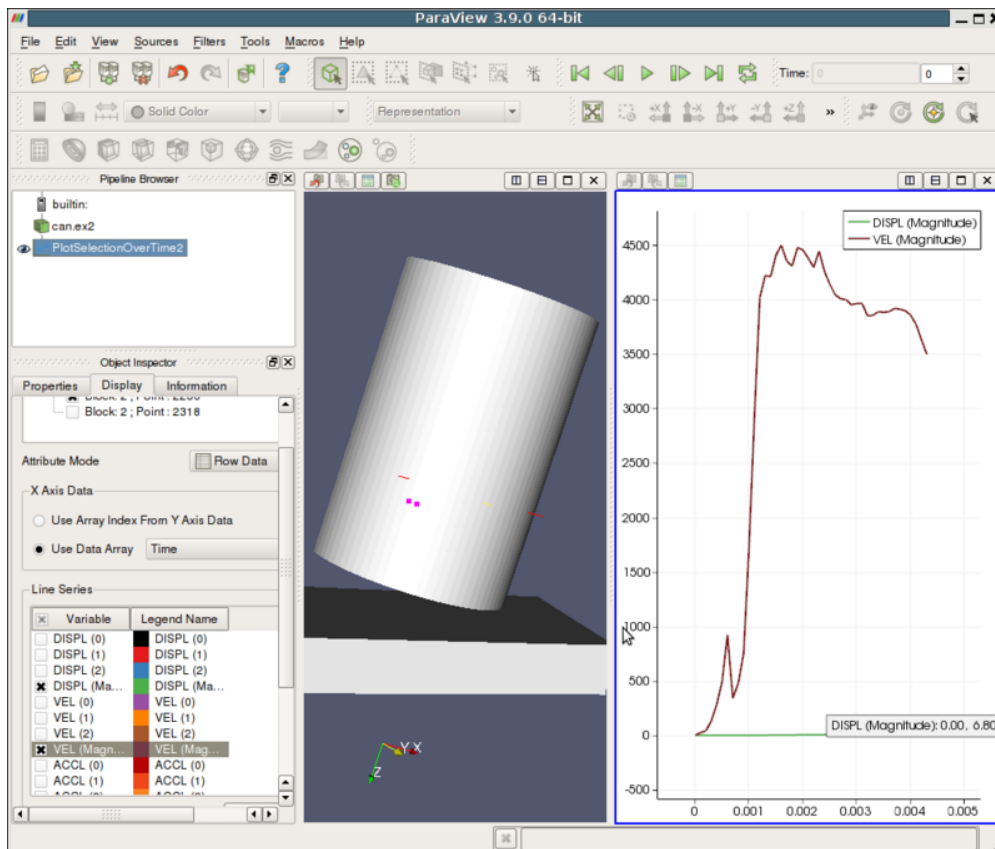
## Extract Selection



Selection makes it possible to highlight and observe regions of interest. Oftentimes once a region of interest has been identified, one would like to apply additional operations on it, such as apply filters to only the selected section of the

data. This can be achieved using the Extract Selection filter. To set the selection to extract, create a selection using any of the methods already described. Then apply the extract selection filter to the source producing the selected data. To copy the active selection to the filter, use the Copy Active Selection button. One can change the active selection at any time and update the filter to use it by using this button. Figure [5] shows the extract selection filter applied after a frustum selection operation. Now, one can treat this as any other data source and apply filters to it, save state, save data etc.

## Plot Selection over Time

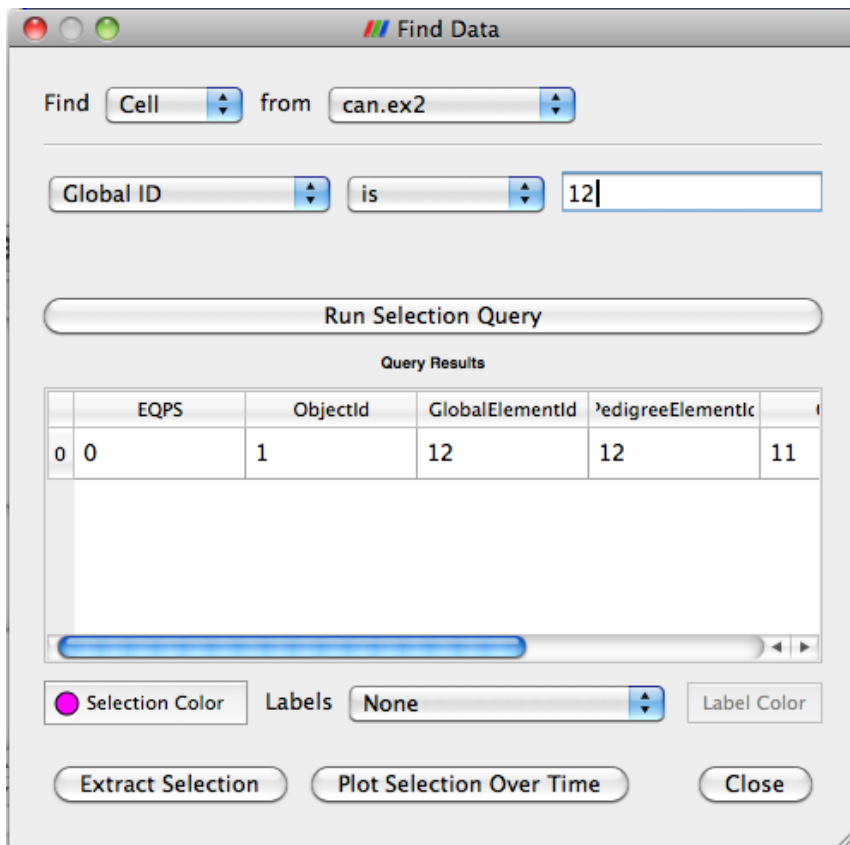


For time varying datasets, you may want to analyze how the data variables change over time for a particular cell or a point. This can be done using the Plot Selection Over Time filter. This filter is similar to the Extract Selection filter, except that it extracts the selection for all time-steps provided by the data source (typically a reader) and accumulates the values for all the cell (or point) attributes over time. Since the selection can comprise of multiple cells or points, the display tab provides the Select Block widget which can be used to select the cell or point to plot, as shown in Figure [6]. Currently only one cell (or point) can be plotted at once in the same xy-plot view. One can create multiple plot views to show multiple plots simultaneously.

# Querying for Data

As previously described, Selection is a mechanism in ParaView for sub-setting and focusing on a particular elements in the dataset. Different views provides different mechanisms for selecting elements, e.g. one can select visible cells or points using the 3D View. Another mechanism for creating selections is by specifying a selection criteria. For example, suppose you want to select all cell where the pressure value is between a certain threshold. In such cases, one can use the **Find Data** dialog. The **Find Data** dialog performs a dual role: not only does it enable specifying the selection criteria but also show details of the selected elements in a spreadsheet. This makes it easier to inspect the selected elements.

To open the **Find Data** dialog, choose **Find Data** from the **Edit** menu.



## When to use Find Data

This feature is useful every time you run into situations where you want to know the cell or the point at which a certain condition happens eg.

- What are the cells at which PRESSURE  $\geq$  12?
- What are the points with TEMP values in the range (12, 133)?
- Locate the cell at ID 122, in Block 2.

This feature provides a convenient way of creating selections based on certain criteria that can then be extracted from the data if needed.



## Using the Find Data dialog

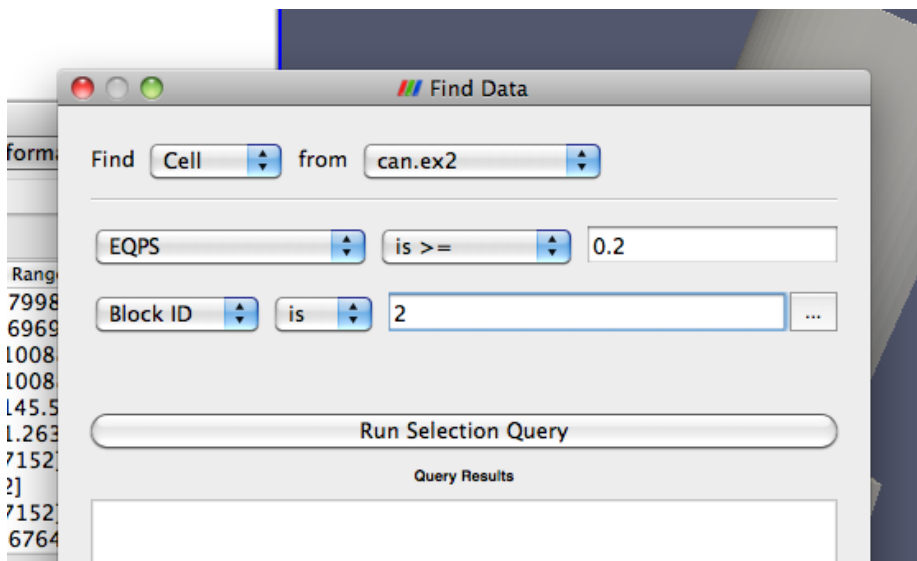
The dialog is designed to be used in two distinct operations:

- Define the selection criteria or query
- Process the selected cells/points e.g. show labels in active 3D view, extract selection etc.

One must define the selection query and then execute the query, using **Run Query** button before being able to inspect or process the selected elements.

## Defining the Query

First decide what type of elements you are interested in selecting i.e. cells or points and from what data-source. This can be done using the following combo-boxes. Note that as you change these, any previous selections/queries will be cleared.



Next, we must define the selection criteria. For defining the criteria, one first choose the term e.g. IDs, Global IDs, attribute array names etc. available for the chosen element type. Next one chooses the operator, and finally the value or values. Based on the operator one or more values may be entered in the value entry box using comma (,) as the separator.

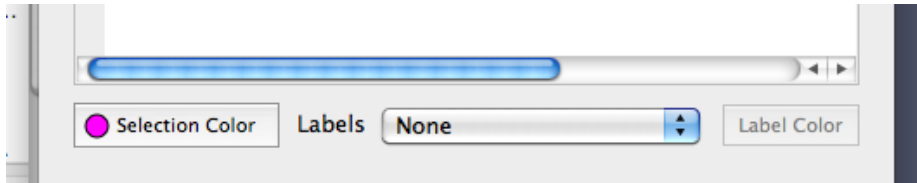
In addition based on the data-type and the nature of the current session, there may additional rows that allow users to qualify the selection using optional parameters such as process number (when running in parallel), or block number (for composite datasets).

Once you have defined your query, hit **Run Query** button to execute the query. If any elements get selected, then they will be shown in the spreadsheet in this dialog. Also, if any of the views are showing the dataset that is selected, then they will highlight the selected elements as well, just like regular view-based selection.



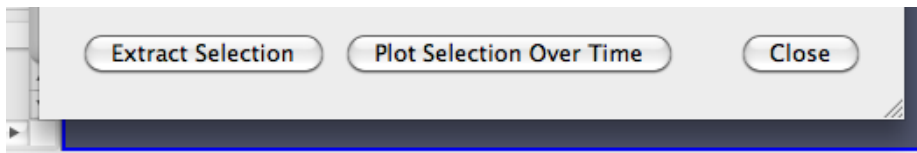
## Displaying the Selection

Once a query is executed, the selected elements will be highlighted in all views where the selected data is visible. If the active view is a 3D view, one can choose whether to show labels for selected elements as well as the color to use for showing the selected elements using the controls on the Find Data dialog itself.



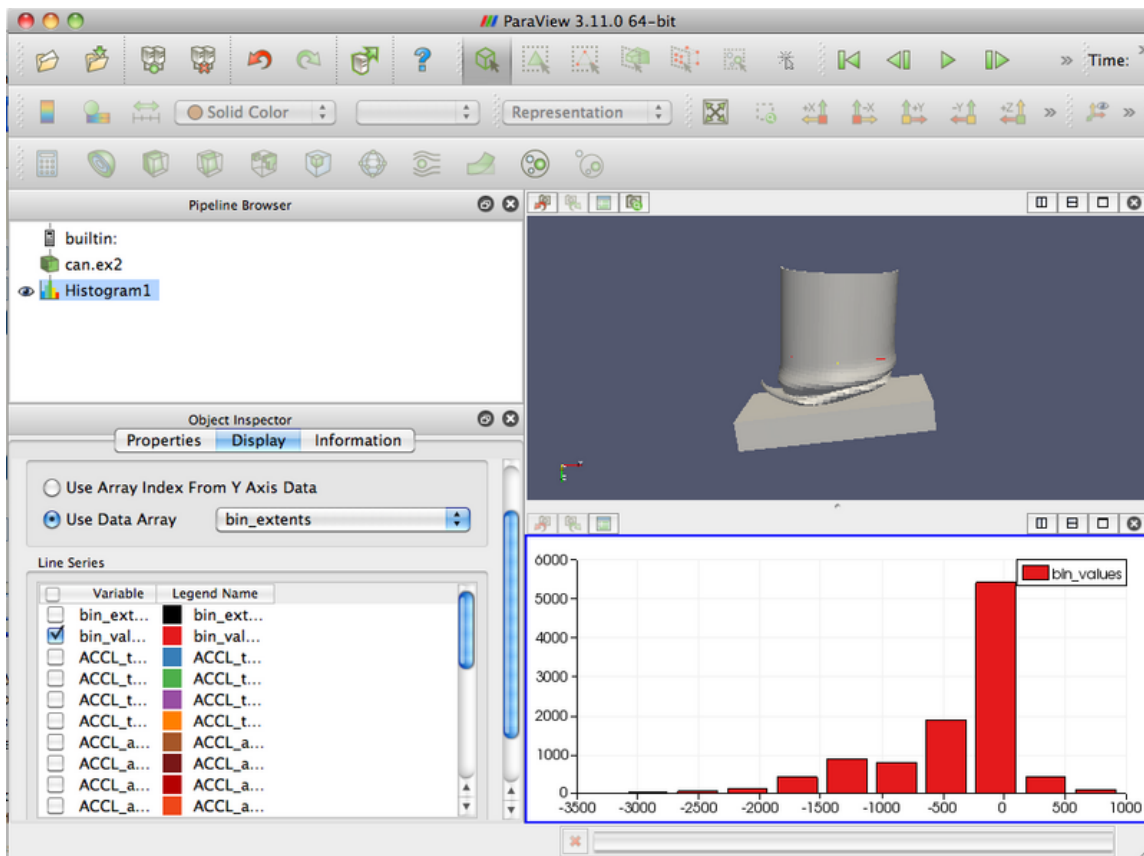
## Extracting the Selection

The results of a query are temporary. They get replaced when a new query is executed or when the user creates a selection using any of the selection mechanisms. Sometimes however, users may want to further analyze the selected elements such as apply more filters to only the selected elements, or plot the change in attributes on the selected elements over time. In that case, one should extract the selection. That creates a new filter that is setup to run the query on its input and produce a dataset matching the selection criteria. Both **Extract Selection** and **Plot Selection Over Time** are filters available through the **Filters** menu. The Find Data dialog provides shortcut buttons to quickly create those filters and set them up with the selection criteria chosen.



## Histogram

The Histogram filter produces output indicating the number of occurrences of each value from a chosen data array. It takes in a `vtkDataObject` but the object must have either point or cell data to operate on. The filter cannot be directly used with a table but the Table to Points filter can be used to convert the table to a polydata for input into the Histogram filter. The bar chart is the default view for the output of the Histogram filter. Other views that can be used are the line chart view, parallel coordinate view, and spreadsheet view. The chart views are useful for observing trends in the data while the spreadsheet view is useful for seeing exact numbers. An example Histogram filter output in a bar chart view is shown below.



Options for the Histogram filter are:

- Which array to process. This can be either a point data array or a cell data array. For arrays with more than 1 component the user can specify which component to compute with respect to. The default is the first component.
- The number of bins to put the results in as well as the range the bins should be divided from.
- An option to average other field information of the same type for each of the bins.
- Which variables are to be displayed in the view. This is under the Display tab of the Object Inspector.

# Plotting and Probing Data

There are multiple ways of probing a data set for point or cell values. The simplest is the Probe Location filter. Additionally, there are a variety of filters to plot data with respect to time, location, or grid object Id.

## Probe Filter

The probe filter can be used to query a data set for point or cell data.

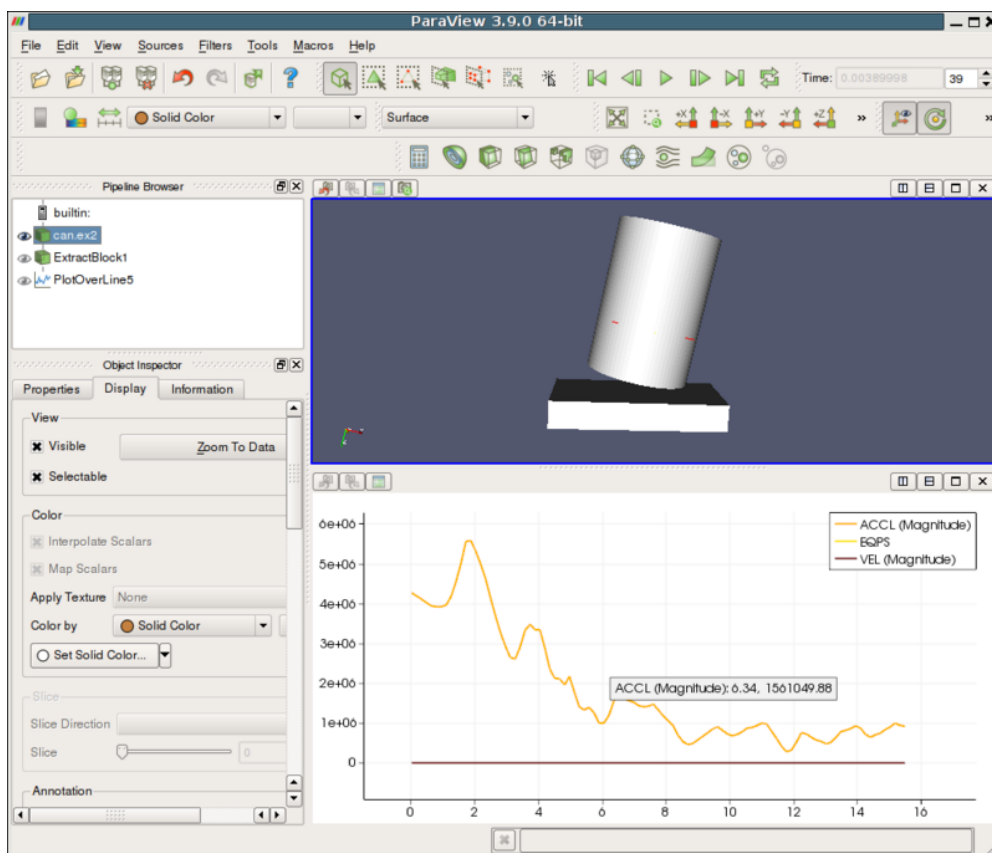
Options for the probe filter are:

- Show point button is used to show where the center of the sphere is for generating the probe points.
- Center on bounds will place the sphere center at the center of the data set's bounds.
- The sphere center can either be set in the properties tab of the object inspector or in the 3d view with the mouse cursor. To choose a new point in the 3d view window, press P and then click on the desired point. Note that the Show Point option must be selected in order to activate the point widget.

The output for this filter is a single cell comprised of a set of points in a vtkPolyData.

## Line Plots

The line plots have similar control functionality for displaying the results. These options are in the Display tab of the Object Inspector and include selecting which point and/or cell data to include in the plot, what should be used as the x axis values (e.g. point/cell index or a specified point or cell data array), and plot preferences such as line thickness and marker style. See the line chart view section for more information on setting the display properties for these filters. An example line plot is shown below.



## Plot Data / Scatter Plot

The Plot Data filter and Scatter Plot filter are very similar in functionality with the Scatter Plot filter being a deprecated version of the Plot Data filter. The main difference is that the Scatter Plot filter's output is a 1D rectilinear grid while the Plot Data filter's output is the same type as the input.

The Plot Data filter plots point or cell data over the entire data set. By default the x axis values are determined with respect to their point or cell index.

## Plot Over Line

The Plot Over Line filter is used to plot point data over a specified straight line. The Plot Over Line filter will not display cell data over the line so the user can use the Cell Data to Point Data filter to convert cell data to point data in order to use this filter for viewing desired cell data. The line geometry can be specified either by setting the points in the Properties tab of the Object Inspector or by pressing P and selecting the beginning and ending points of the line. Note that the Show Line option must be selected to activate the line widget. The resolution option is used to specify at how many evenly spaced points on the line to query the data set for information the point data to plot with respect to.

## Plot on Sorted Lines

The Plot on Sorted Lines filter is used when polydata with line cell types have been extracted from a data set and the user wants to see how point data varies over the line. This filter orders the lines by their connectivity instead of their point index. The output of the filter is a multiblock with a block for each contiguous line.

## Plot on Intersection Curves

The Plot on Intersection Curves filter is used to plot point data where the data set intersects the given polygonal slice object. The results is a multiblock of polydatas where each block represents a contiguous set of cells. Note that the polydatas will only have 1D cell types. Thus if the slice type object intersection with the data set has 2D geometry the filter will take the boundary of the slice type object in order to reduce down to 1D geometry. If the slice type is a sphere which is wholly contained in a volumetric data set then the filter parameters are invalid and no lines will be output.

## Plot Selection Over Time

The Plot Selection Over Time filter can be used to visualize the variation of point or cell data from a selection with respect to time. The selection should be made on the data set of the filter that precedes the Plot Selection Over Time filter in the pipeline in order to ensure that the proper points or cells are selected. Note that information for only a single point or cell can be plotted at a time.

## Plot Global Variables Over Time

The Plot Global Variables Over Time filter plots field data that is defined over time. Filters must be set up explicitly to provide the proper information as this will not work for field data in general. As an example see the Exodus reader in ParaView.

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# Saving Data

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## Saving Data

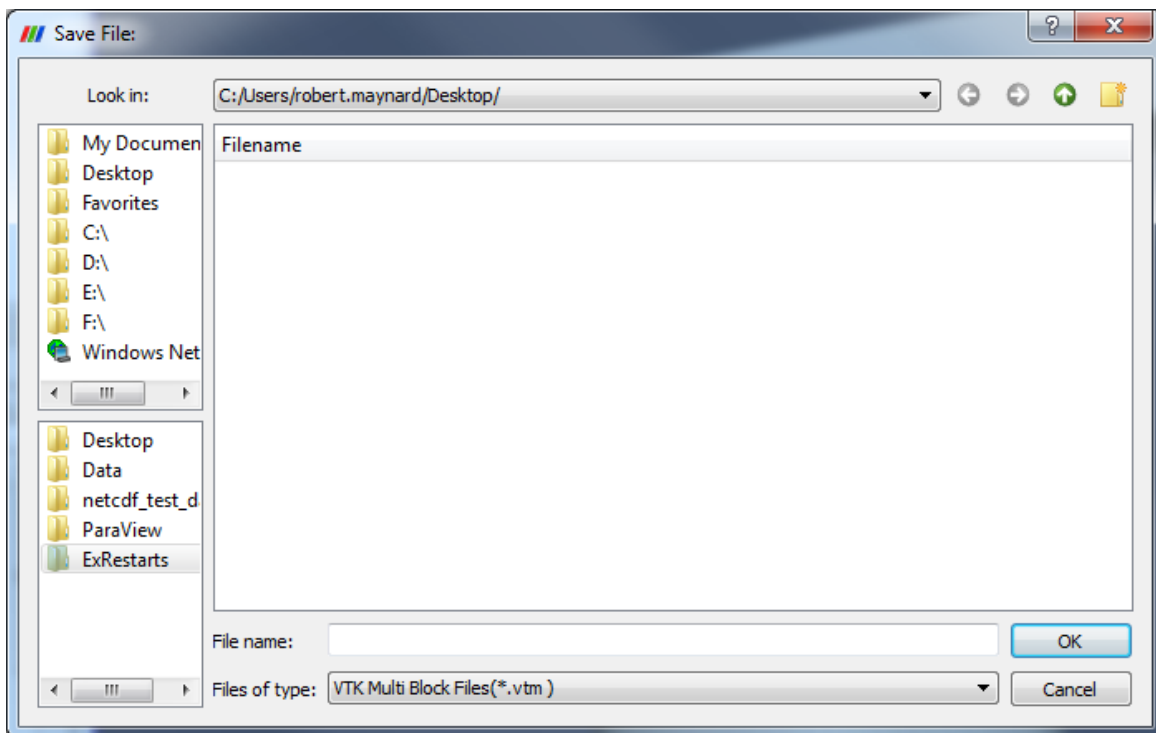
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### Saving Data

Once you have created a visualization of your data in ParaView, you can save the resulting work as raw data, image, movie, and geometry.

#### Save raw data

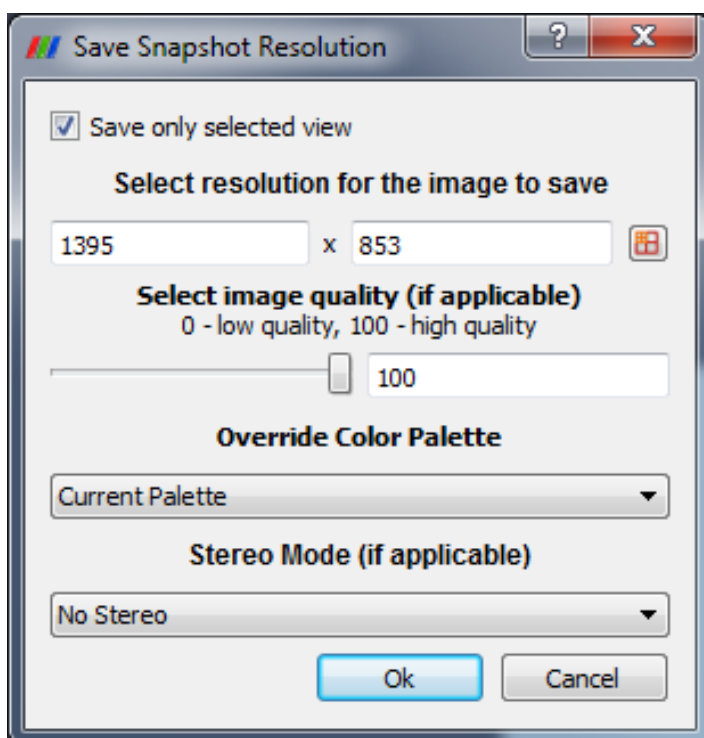
Any object in the ParaView pipeline browser can be saved to a data file by selecting **Save Data** from the **File** menu. The available file types will change based on the data set type of the current data set. The file formats in which ParaView can save data are listed on List of Writers



#### Save screenshots

ParaView allows the user to save either the active view or all the views. The resulting image will be saved on the client machine, even when running with a remote server. The dialog allows you to control the following:

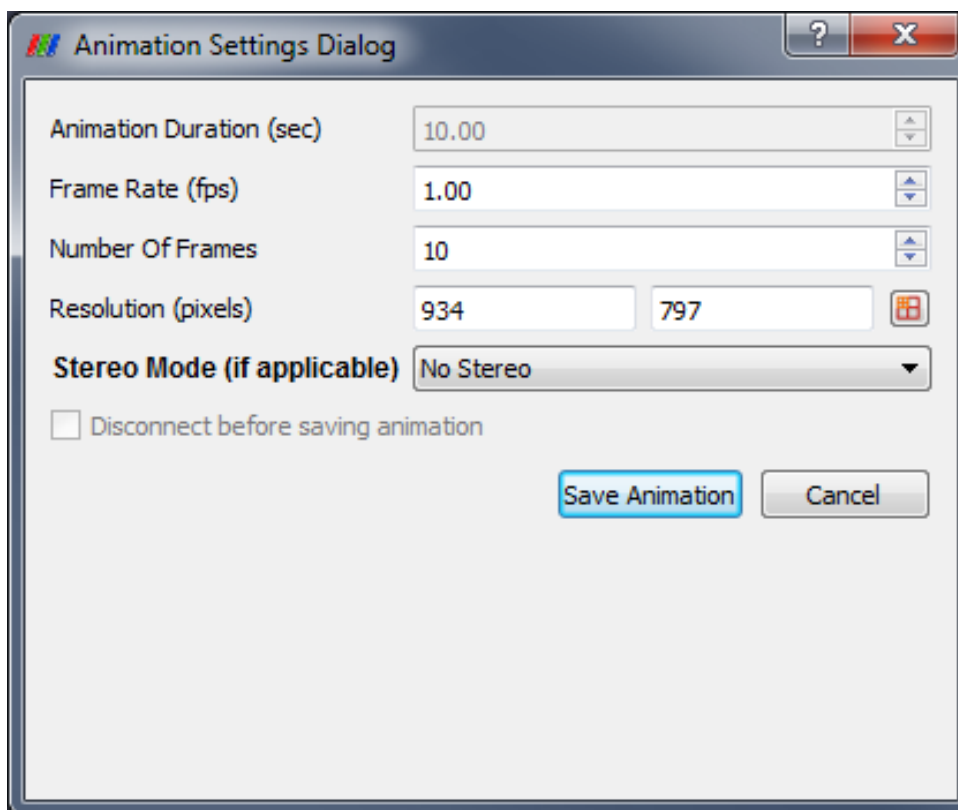
- Image size
- Aspect Ratio of the Image
- Image Quality
- Color Palette
- Stereo Mode



## Save Animation

Once you have created an animation of your data, you can save the animation to disk either as a series of images (one per animation frame) or as a movie file. The animation will contain all the visible views.

To do this, select **Save Animation** from the **File** menu. The **Animation Settings Dialog** then appears, which lets you set properties for the recorded animation.



Once you press the Save Animation button, a save file dialog box will allow you to choose where to save your image or movie file(s). Enter a file name, select an image file type (.jpg, .tif, or .png) or a movie file type (.avi).

Once you choose a file name, the animation will play from beginning to end, and the animation frames generated will be used to create the selected type of image or movie file(s). While the image is playing the render window in ParaView will not update with the correct frame.

If you are connected to a remote server then the Disconnect before saving animation checkbox will be enabled. If you select this before saving the animation, then the ParaView client will disconnect from the server immediately, and the server will continue generating and saving images until the animation completes. When it is finished recording the animation, the server will shut down.

## Save geometries

In addition to saving images of each time step in your animation, you may also wish to save the geometry itself. Select **Save Geometry** from the **File** menu to do this. This will cause a file navigation dialog box to be displayed. Navigate to the location where you wish to save the geometry, and enter a file name. You must save your data using ParaView's .pvd file format.

Unlike an animation, save geometry will only save the visible geometry of the active view for each time step. The resulting .pvd file will contain a pointer to each of these files, which are saved in a folder with the same as the .pvd file.

You can later reload the .pvd file into ParaView as a time varying data set. If multiple data sets were displayed while the animation was running, they will be grouped together as a multi-block data set for each time step. If you then want to operate on the parts individually, run the **Extract Blocks** filter to select the appropriate block(s).

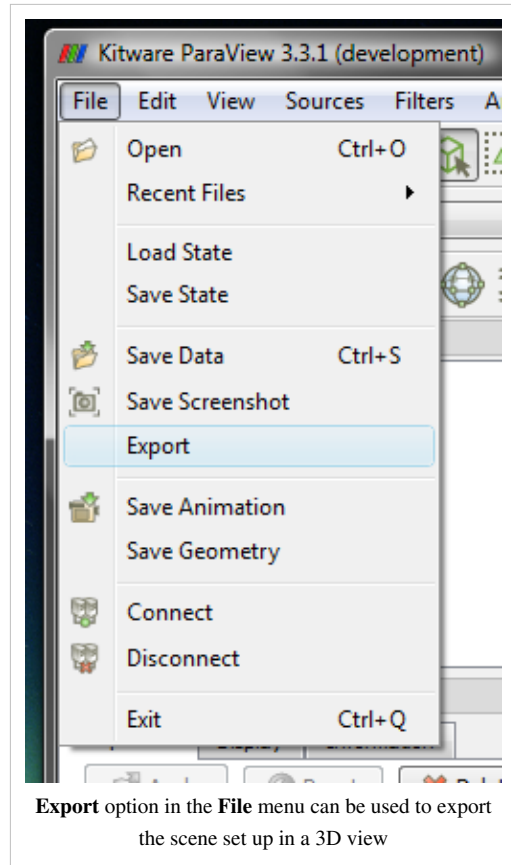
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# Exporting Scenes

ParaView provides functionality to export any scene set up with polygonal data (i.e without volume rendering). Currently X3D <sup>[1]</sup> (ASCII as well as binary), VRML (Virtual Reality Modeling Language) <sup>[2]</sup>, and POV-Ray <sup>[3]</sup> are supported. To export a scene, set up the scene in a 3D view. Only one view can be exported at a time. With the view to be exported active, choose **File | Export**.

The file-open dialog will list the available types. The type is determined based on the extensions of the file written out:

- \*.vrml -- VRML <sup>[4]</sup>
- \*.x3d -- X3D ASCII <sup>[5]</sup>
- \*.x3db -- X3D Binary <sup>[5]</sup>
- \*.pov -- POV-Ray <sup>[6]</sup>



**Export** option in the **File** menu can be used to export the scene set up in a 3D view

ParaView: [Welcome | Site Map <sup>[7]</sup>]

## References

- [1] <http://www.web3d.org/x3d/>
- [2] <http://www.web3d.org/x3d/vrml>
- [3] <http://www.povray.org/>
- [4] <http://www.vtk.org/doc/nightly/html/classvtkVRMLExporter.html>
- [5] <http://www.vtk.org/doc/nightly/html/classvtkX3DExporter.html>
- [6] <http://www.vtk.org/doc/nightly/html/classvtkPOVExporter.html>
- [7] <http://www.paraview.org/Wiki/Category:ParaView>



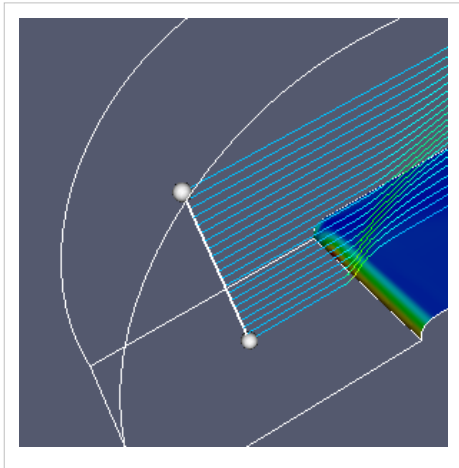
# 3D Widgets

## Manipulating data in the 3D view

### 3D Widgets

In addition to being controlled manually through entry boxes, sliders, etc., parameters of some of the filters and sources in ParaView can be changed interactively by manipulating 3D widgets in a 3D view. Often the 3D widgets are used to set the parameters approximately, and then the manual controls are used for fine-tuning these values. In the manual controls for each 3D widget, there is a check box for toggling whether the 3D widget is drawn in the scene; the label for the check box depends on the type of 3D widget being used. The following 3D widgets are supported in ParaView.

### Line Widget



The line widget is used to set the orientation and the position of a line. It is used in both the **Stream Tracer** and **Elevation** filters. The position of the line can be changed by clicking on any point on the line except the endpoints and dragging. To position the widget accurately, the user may need to change the camera position as well. Holding the Shift key while interacting will restrict the motion of the line widget to one of the X, Y, or Z planes. (The plane chosen is the one most closely aligned with the direction of the initial mouse movement.) To move one of the endpoints, simply use one of the point widgets on each end of the line. These are marked by spheres which become red when clicked. You can also reposition the endpoints by pressing the “P” key; the endpoint nearest to the mouse cursor will be placed at the position on the data set surface beneath the mouse position. Left-clicking while the cursor is over the line and dragging will reposition the entire line. Doing the same with the right mouse button causes the line to resize. Upward mouse motion increases the length of the line; downward motion decreases it.

|   |                                  |                                  |                                  |
|---|----------------------------------|----------------------------------|----------------------------------|
| <input checked="" type="checkbox"/> Show Line |                                  |                                  |                                  |
| Point1  | <input type="text" value="-10"/> | <input type="text" value="-10"/> | <input type="text" value="-10"/> |
| Point2  | <input type="text" value="10"/>  | <input type="text" value="10"/>  | <input type="text" value="10"/>  |
| <input type="text" value="X Axis"/>           |                                  |                                  |                                  |
| <input type="text" value="Y Axis"/>           |                                  |                                  |                                  |
| <input type="text" value="Z Axis"/>           |                                  |                                  |                                  |
| Resolution                                    | <input type="text" value="100"/> |                                  |                                  |

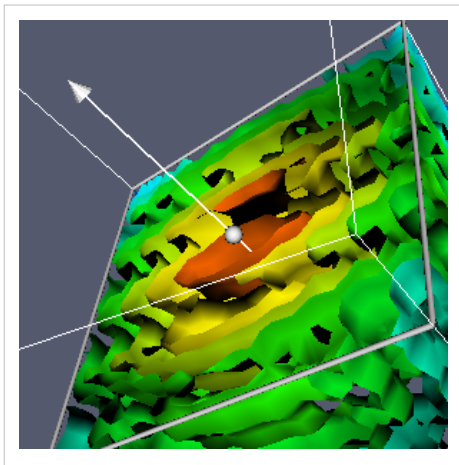
Figure 10. Line widget user interface

The **Show Line** check box toggles the visibility of the line in the 3D view.

The controls shown above can be used to precisely set the endpoint coordinates and resolution of the line. The **X Axis**, **Y Axis**, and **Z Axis** buttons cause the line to be along the selected axis and pass through the center of the bounds of the data set.

Depending on the source or filter using this widget, the **Resolution** spin box may not be displayed. The value of the **Resolution** spin box determines the number of segments composing the line.

## Plane Widget



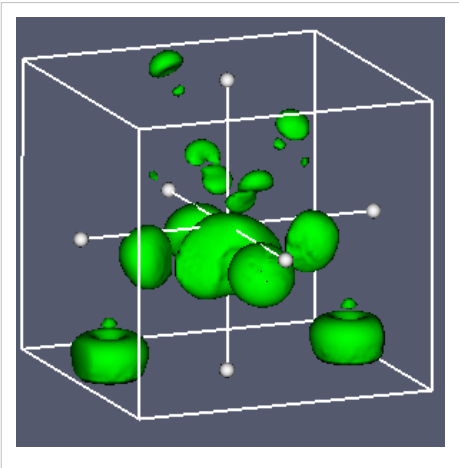
The plane widget is used in clipping and cutting. The plane can be moved parallel to its normal by left-clicking on any point on the plane except the line center and dragging. Right-clicking on the plane (except on the normal line center) and dragging scales the plane widget. Upward mouse motion increases the size of the plane; downward motion decreases it. The plane normal can be changed by manipulating one of the point widgets (displayed as cones that become red when clicked) at each end of the normal vector.

Shown below, the standard user interface for this widget provides entry boxes for setting the center position (**Origin**) and normal direction of the plane as well as toggling the plane widget's visibility (using the **Show Plane** check box). Buttons are provided for positioning the plane at the center of the bounding box of the data set (**Center on Bounds**) and for aligning the plane's normal with the normal of the camera (**Camera Normal**), the X axis, the Y axis, or the Z axis (**X Normal**, **Y Normal**, and **Z Normal**, respectively). If the bounds of the data set being operated on change, then you can use the **Reset Bounds** button to cause the bounding box for the plane widget to match the new bounds of the data set and reposition the origin of the plane at the center of the new bounds. (Using only the **Center on Bounds** button in this case would move the origin to the center of the new bounds, but the bounds of the widget would not be updated.)

|  |                                |   |
|--|--------------------------------|---|
| <input checked="" type="checkbox"/> Show Plane |                                |   |
| Origin   | <input type="text" value="0"/> | <input type="text" value="0"/>                  |
| Normal   | <input type="text" value="1"/> | <input type="text" value="0"/>                  |
| <input type="button" value="X Normal"/>        |                                | <input type="button" value="Reset Bounds"/>     |
| <input type="button" value="Y Normal"/>        |                                |   |
| <input type="button" value="Z Normal"/>        |                                |   |
| <input type="button" value="Camera Normal"/>   |                                | <input type="button" value="Center on Bounds"/> |

Figure 11. Plane widget user interface

Box Widget



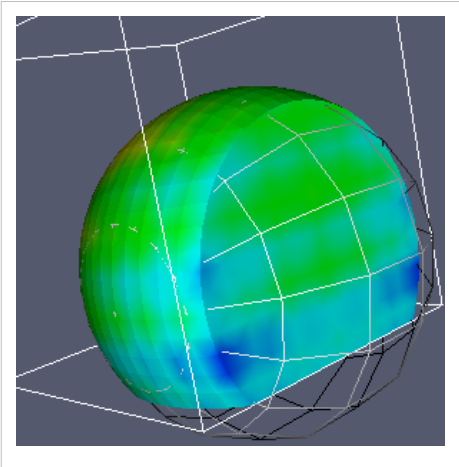
The box widget is used for clipping and transforming data sets. Each face of the box can be positioned by moving the handle (sphere) on that face. Moving the handle at the center of the box causes the whole box to be moved. (This can also be achieved by holding the Shift key while interacting.) The box can be rotated by clicking and dragging with the left mouse button on a face (not at the handle) of the box. Clicking and dragging inside the box with the right mouse button uniformly scales the box. Dragging upward increases the box's size; downward motion decreases it.

Traditional user interface controls, shown below, are also provided if more precise control over the parameters of the box is needed. These controls provide the user with entry boxes and thumb wheels or sliders to specify translation, scaling, and orientation in three dimensions.

|  |                                |                                |                                |
|--|--------------------------------|--------------------------------|--------------------------------|
| Translate:                                     | <input type="text" value="0"/> | <input type="text" value="0"/> | <input type="text" value="0"/> |
| Scale:   | <input type="text" value="1"/> | <input type="text" value="1"/> | <input type="text" value="1"/> |
| Orientation:                                   | <input type="text" value="0"/> | <input type="text" value="0"/> | <input type="text" value="0"/> |
| <input checked="" type="checkbox"/> Visibility |                                |                                |                                |

Figure 12. Box widget user interface

Sphere Widget



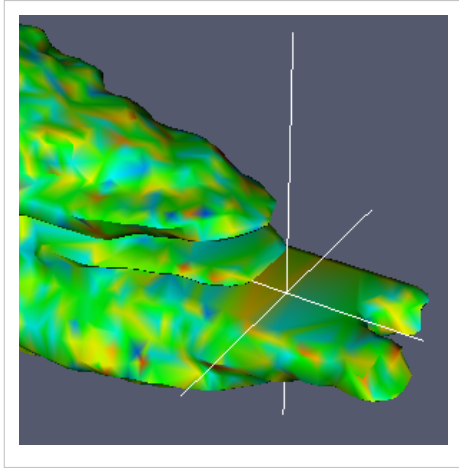
The sphere widget is used in clipping and cutting. The sphere can be moved by left-clicking on any point on the sphere and dragging. The radius of the sphere is manipulated by right-clicking on any point on the sphere and dragging. Upward mouse motion increases the sphere's radius; downward motion decreases it.

As shown below, the center and radius can also be set manually from the entry boxes on the user interface. There is also a button to position the sphere at the center of the bounding box of the current data set.

|  |                                      |                                      |  |
|--|--------------------------------------|--------------------------------------|--|
|  | x                                    | y                                    | z                                      |
| Center   | <input type="text" value="-4.8784"/> | <input type="text" value="-4.7546"/> | <input type="text" value="-0.020147"/> |
| Radius   | <input type="text" value="10"/>      |                                      |  |
| <input type="button" value="Set Sphere Center to Center of Bounds"/> |                                      |                                      |  |
| <input checked="" type="checkbox"/> Visibility                       |                                      |                                      |  |

Figure 13. Sphere widget user interface

## Point Widget

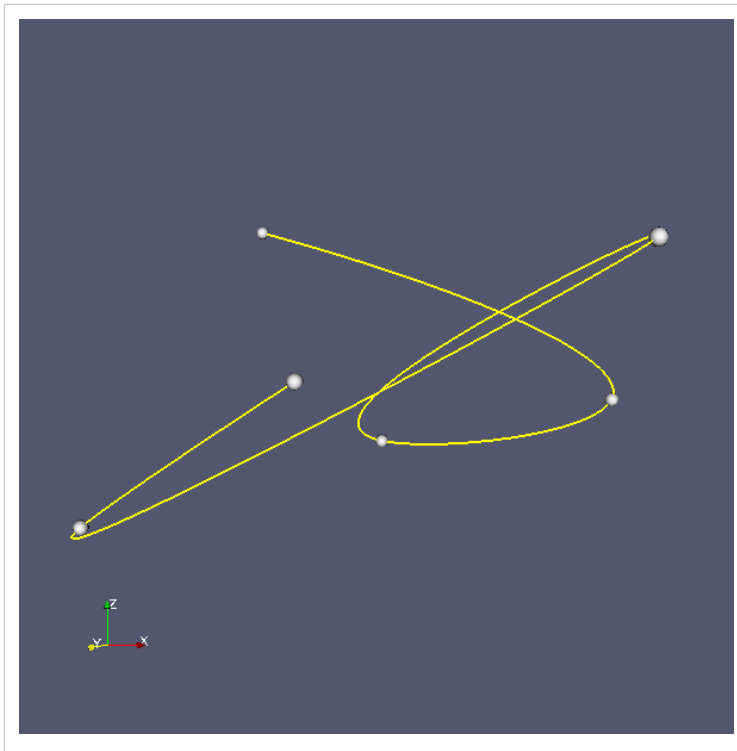


The point widget is used to set the position of a point or the center of a point cloud. It is used by both the **Stream Tracer**, **Probe Location** and **Probe Location over Time** filters. The position of the point can be changed by left-clicking anywhere on it and dragging. Right-clicking and dragging anywhere on the widget changes the size of the point widget in the scene. To position the widget accurately, the user may need to change the camera position as well. Holding the Shift key while interacting will restrict the motion of the point to one of the x, y or z planes. (The plane chosen is the one that is most closely aligned with the direction of the initial mouse movement.)

As shown below, entry boxes allow the user to exactly specify the coordinates of the point, and a button is provided to position the point at the center of the bounds of the current data set. If the point widget is being used to position a point cloud instead of a single point, entry boxes are also provided to specify the radius of the point cloud and the number of points the cloud contains.

Figure 14. Point widget user interface

## Spline Widget



The spline widget is used to define a path through 3D space. It is used by the **Spline Source** source and in the Camera animation dialog. The widget consists of a set of control points, shown as spheres in the 3D scene that can be clicked on with the mouse and dragged perpendicular to the viewing direction. The points are ordered and the path through them defines a smoothly varying path through 3D space.

As shown below, the text control box for the spline widget allows you to add or delete control points and specify their locations exactly. You can hide or show the widget and can choose to close the spline to create a loop, which adds a path segment from the last control point back to the first.

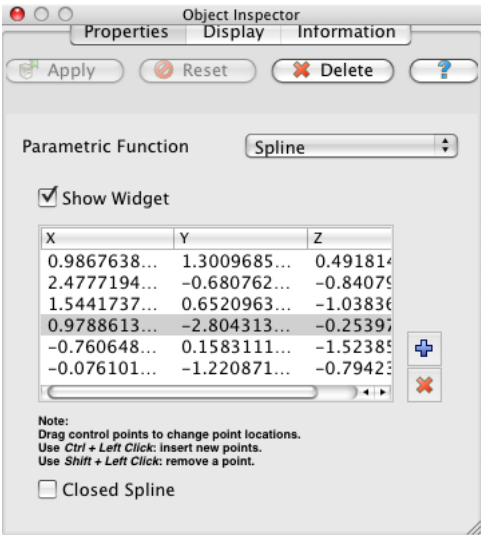


Figure 15. Spline widget user interface

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# Annotation

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## Annotation

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### Annotation

In ParaView, there are several ways to annotate the data to better understand or explain it. Several of the annotations can be interactively placed in the 3D scene. The parameters of the annotations are controlled using traditional user interface elements. Some types of annotation are discussed in other parts of this book. See the section on Selection for information about labeling selected points or cells in a data set.

### Scalar Bar

The most straightforward way to display a scalar bar (or color legend) in the active 3D view is to use the color legend visibility button  on the **Active Variables Control** toolbar. When the data set selected in the **Pipeline Browser** is being colored by a variable (i.e., something other than **Solid Color** is selected in the **Color by** menu on the **Display** tab), this button is active. Clicking it toggles the visibility (in the selected 3D view) of a scalar bar showing the mapping from data values to colors for that variable.

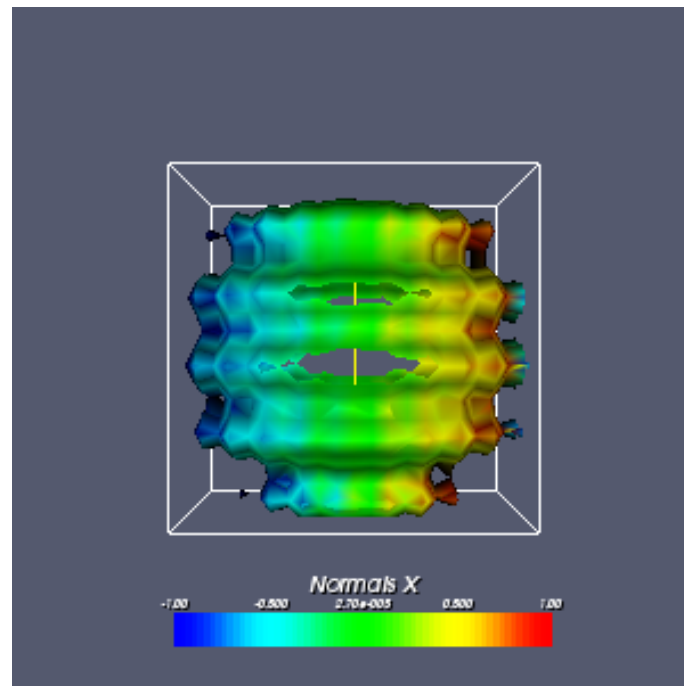



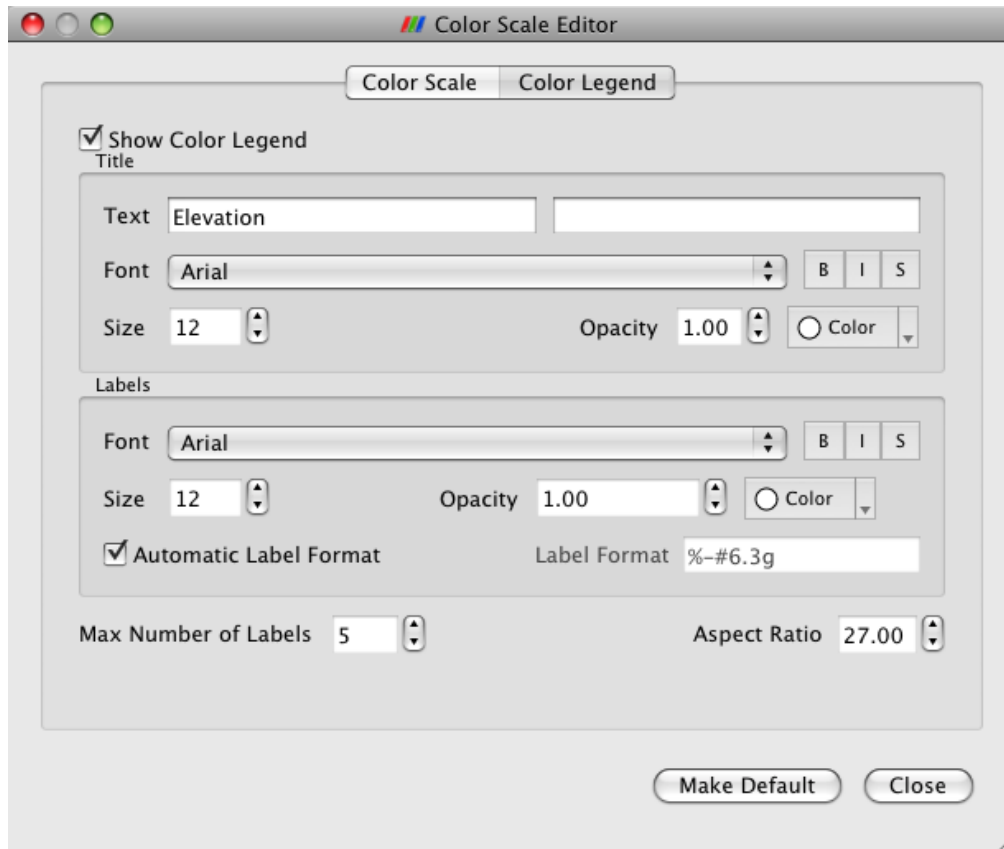
Figure 4. Scalar Bar displaying X-component of Normals

Clicking the **Edit Color Map** button to the right of the  **Color Legend** button brings up the **Color Scale Editor** dialog. As described in the Displaying Data <sup>[1]</sup> chapter you have precise control over the mapping between data values and visible colors on the **Color Scale** tab of that dialog.

On the **Color Legend** tab, there are several different parameters that allow you to control the appearance of the color legend itself. First is the **Show Color Legend** check box which toggles the visibility of the scalar bar (color legend) in the 3D view. This has the same effect as using the color legend visibility button. Below the **Show Color Legend** check button are the controls for displaying the title and labels on the scalar bar. In the **Title** section, the **Text** entry

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box specifies the label that appears above the scalar bar. It defaults to the name of the array used for coloring the data set. If the current data set is being colored by a vector array, the value of the second entry box defaults to specifying how the color is determined from the vector (i.e., **X**, **Y**, **Z**, or **Magnitude**). The entry box labeled **Labels** contains formatting text specifying the form of the scalar bar labels (numbers). The format specification used is the same as that used by the `printf` function in C++.



**Figure 5. The color legend tab of the Color Scale Editor dialog**

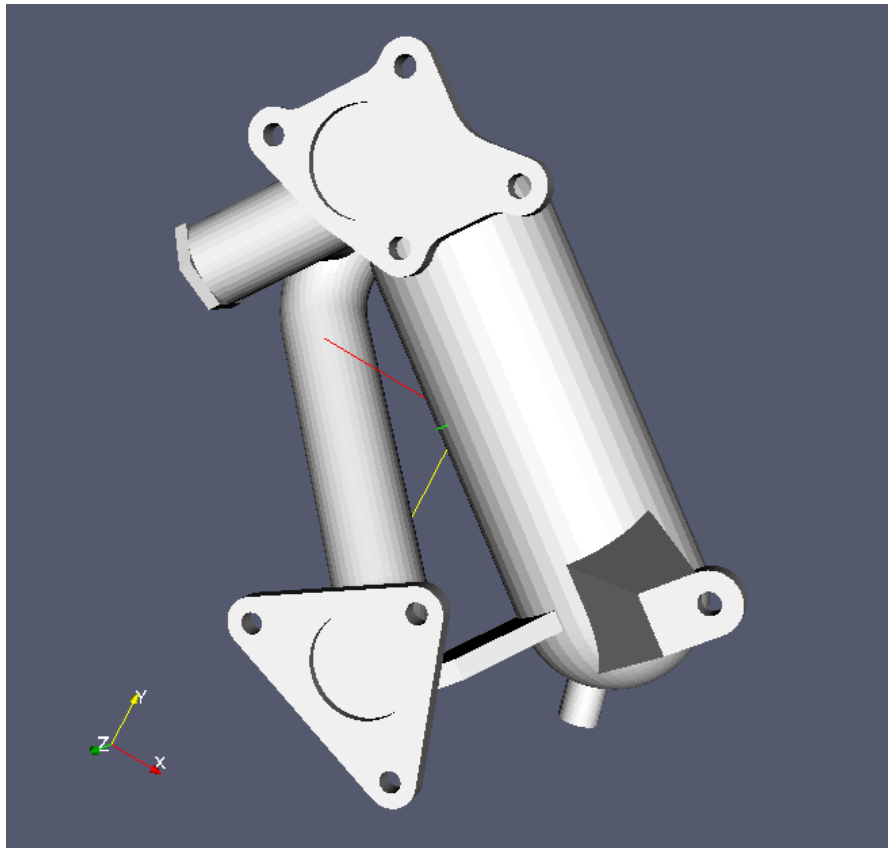
Below each of the **Title** and **Labels** entry boxes are controls for determining how the title and labels will be drawn in the display area. The leftmost control is a menu for choosing the font; the available fonts are **Ariel** (the default), **Courier**, and **Times**. Next are three formatting attribute buttons controlling whether the text is boldfaced, italicized, or shadowed, respectively. The next interface control is a spin box for controlling the text's opacity. It ranges from 0 (transparent) to 1 (opaque).

At the bottom of this tab is a **Number of Labels** spin box. This determines how many scalar-value labels will be shown alongside the scalar bar. To the right of that is an "Aspect Ratio" spin box which allows you to make the scalar bar relatively thinner or thicker.

When the scalar bar is displayed in the 3D scene, it can be positioned and resized interactively, similar to interacting with 3D widgets. Clicking and dragging the scalar bar with the left mouse button repositions it in the display area. If the scalar bar begins to move beyond the left or right side of the display area, it is reoriented vertically. It is reoriented in a horizontal direction if it is moving off-screen at the top or bottom of the display area. The scalar bar can be resized by left-clicking and dragging any of its sides or corners.

## Orientation Axes

When interacting with data in ParaView, it can be difficult to determine how the data set is oriented in 3D. To remedy this problem, a marker showing labeled 3D axes (i.e., orientation axes) can be displayed. The orientation of the axes matches that of the data, and the axes reorient themselves as the camera is rotated about the 3D scene.



**Figure 5. 3D Orientation Axes**

The user interface controls for the orientation axes are located in the **Annotation** section of the **View Settings** dialog (*Edit menu, View Settings*) **for the 3D view**. The **check box beside the 'Orientation Axes** label toggles the visibility of the orientation axes. The **Interactive** check box controls whether the orientation axes can be repositioned and resized through mouse interaction, similar to interacting with 3D widgets. Left-clicking and dragging the orientation axes repositions them in the 3D view. When the orientation axes are in interactive mode, a bounding rectangle (outline) is displayed around the axes when the mouse moves over them. Left-clicking and dragging on the edges of this outline resizes the orientation axes. From the **Set Outline Color** button (enabled when **Interactive** is checked), you can change the color used for displaying the bounding rectangle. This is useful for making the outline more visible if its current color is similar to the color of the background or any object behind the outline (if the orientation axes are placed in front of a data set). The orientation axes are always drawn in front of any data set occupying the same portion of the 3D view. The **Axis Label Color** button allows you to change the color of the labels for the three axes. The reasons for changing the color of the axis labels are similar to those for changing the outline color.



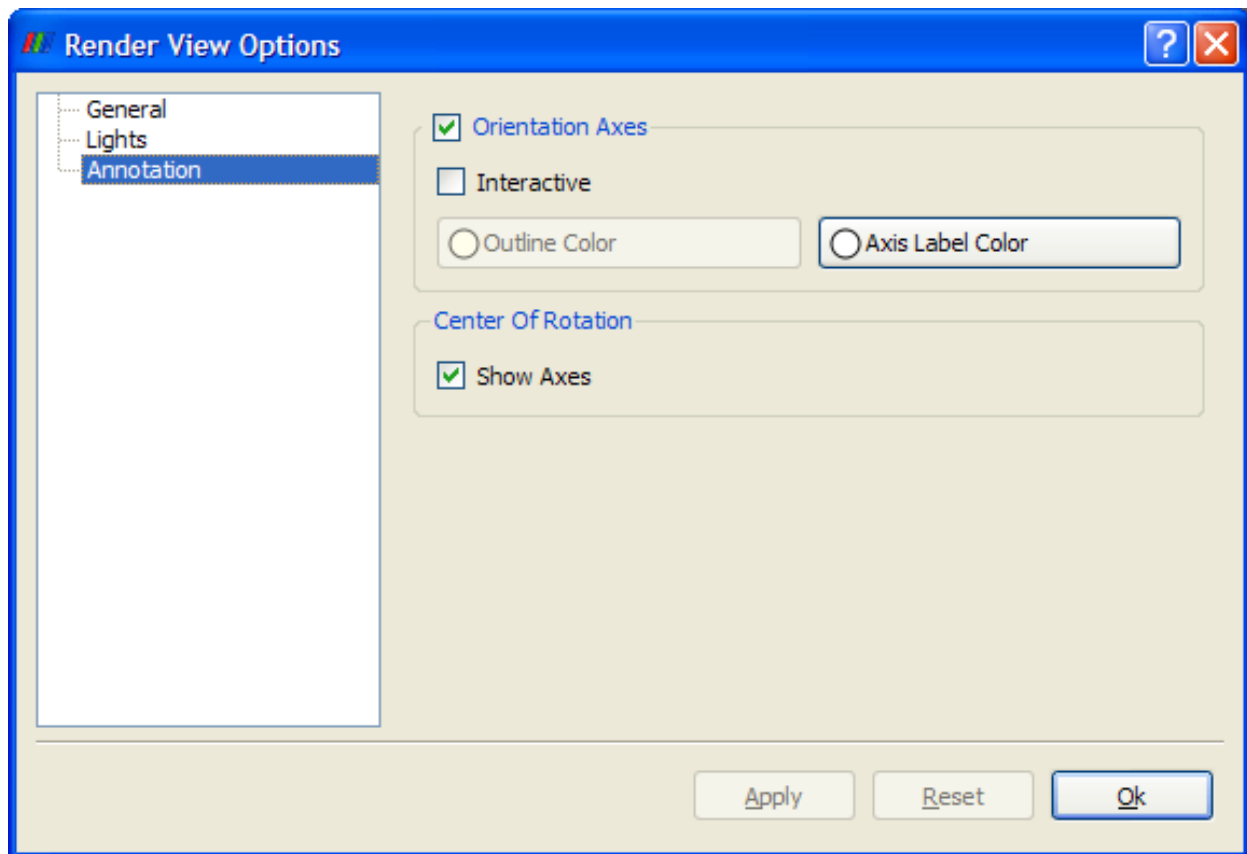


Figure 6. User interface controls for orientation axes (in 3D view's view settings dialog)

## Center of Rotation

On the same dialog box you also have control over whether or not the center of rotation should be annotated. This annotation demarcates the location in space about which the camera rotates. ParaView uses the center of the initial data set loaded into the pipeline as the default value. You can control this placement and display of the center of rotation display via the **Center Axes Control toolbar** (*View menu, Toolbars, 'Center Axes Control'*). If you need to specify an exact coordinate you may do so via the **Adjust Camera** dialog exposed by the rightmost button on the left set of buttons above the 3D view.

## Text Display

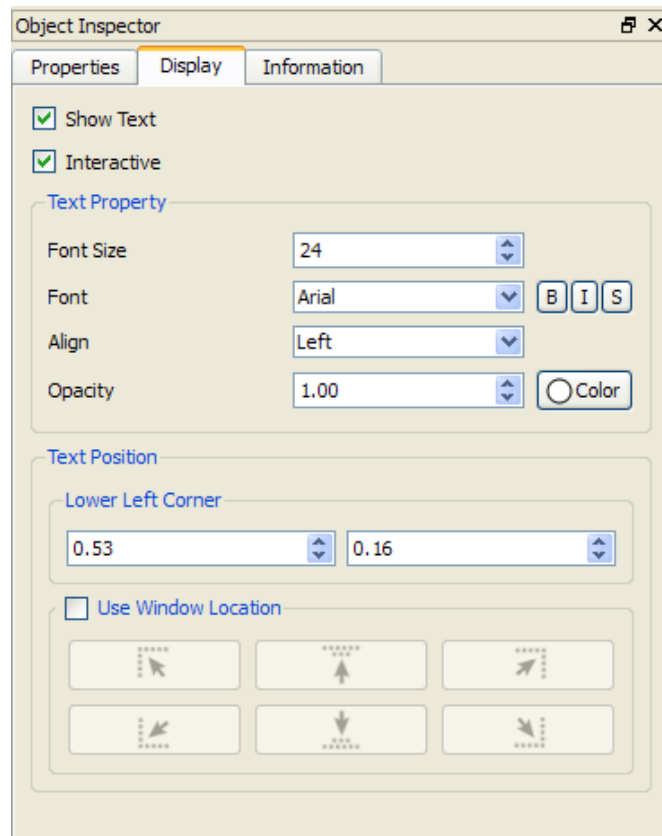
Often users want to display additional text in the 3D view (e.g., when generating a still image or an animation). In ParaView, there are a variety of ways to do this, all through the use of sources and filters. The **Text** source displays 2D text in the 3D scene, the **3D Text** source does the same for 3D text, and the **Annotate Time** filter shows the current time using a text annotation in the 3D view.

### Text Source

The **Text** source allows you display arbitrary 2D text on top of a 3D view. It is available from the **Sources** menu. On its **Properties** tab, enter whatever text you wish to be displayed in the **Text** entry area, and press the **Apply** button.

Once the text has been added to the 3D view, you can reposition it within the view by left-clicking and dragging the text. (A bounding box will be displayed around the text.) By default, the text is displayed in the lower left corner of the 3D view.

The **Display** tab in the **Object Inspector** for the **Text** source is different than for most other sources and filters displayed in the 3D view. Specific controls are given for positioning the text and setting various font properties.



**Figure 7. Display tab for the Text source**

The first check box, **Show Text**, toggles the visibility of the text in the 3D view. This corresponds to the eye icon in the **Pipeline Browser** for this source. The **Interactive** check box below it determines whether the text may be interactively repositioned by clicking and dragging in the 3D view.

The **Text Property** section of the **Display** tab allows you to specify various parameters relating to the font used for displaying the text. The **Font Size** spin box determines how large the letters in the text will be. The **Font** menu determines which font type will be used: **Arial**, **Courier**, or **Times**. The three buttons beside this menu indicate whether the text will be drawn boldfaced, italicized, and/or using shadows. The **Align** menu specifies whether the text will be left-, center-, or right-aligned within its bounds box. The **Opacity** spin box determines how opaque the text appears. An opacity value of 0 corresponds to completely transparent text, while an opacity value of 1 means completely opaque text. The **Color** button beside the Opacity spin box allows you to choose a color for the text.

The **Text Position** of the **Display** tab is used for specifying exactly where in the 3D view the text appears. If **Use Window Location** is unchecked, then the **Lower Left Corner** section is active. The two spin boxes provided determine the X and Y coordinates of the lower left corner of the bounding box of the text, specified in normalized coordinates (starting from the lower left corner of the view).

If **Use Window Location** is checked, then the six buttons in that section of the interface become available. They allow you to choose various locations within the view for anchoring the text. Using the top row of buttons, you may place the text in the upper left corner, in the top center of the view, or in the upper right corner, respectively. The bottom row of button behaves similarly for the bottom portion of the 3D view.

### 3D Text Source

Whereas the **Text** source (described above) positions text in 2D, overlaying the 3D scene, the **3D Text** source places text *in* the 3D scene. The text is affected by rotation, panning, and zooming as is any other object in the 3D scene.

You can control the placement of the 3D Text in two ways. First, the **Display** tab for all sources and filters shown in a 3D view have controls to position, translate, rotate and scale the object via textual controls. It is often easier however to use a the **Transform** filter for the same purpose. The transform filter allows you to drag, rotate and scale the object via either the 3D widget in the scene or via the text controls on its Properties tab.

### Annotate Time

The **Annotate Time** source and filter are useful for labeling a time-varying data set or animation with ParaView's current time in a 3D view. The distinction between the two is somewhat abstract. VTK's temporal support works by have the pipeline request data for a particular time. The text source displays exactly this. VTK's sources and filters are allowed to produce data at different (usually nearby producing data that changes over a step function in time) time. The annotate time filter shows the data produces by an given filter.

In either case, the annotated time value is drawn as a 2D label, similar to the output of the **Text** source. The **Properties** tab for this filter provides a single text box labeled **Format**. From this text box, you may specify a format string (printf style) indicating how the label will appear in the scene. As with the output of the **Text** source, this label may be interactively repositioned in the scene. The **Display** tab for this source is the same as the one for the **Text** source; it is described earlier in this section.

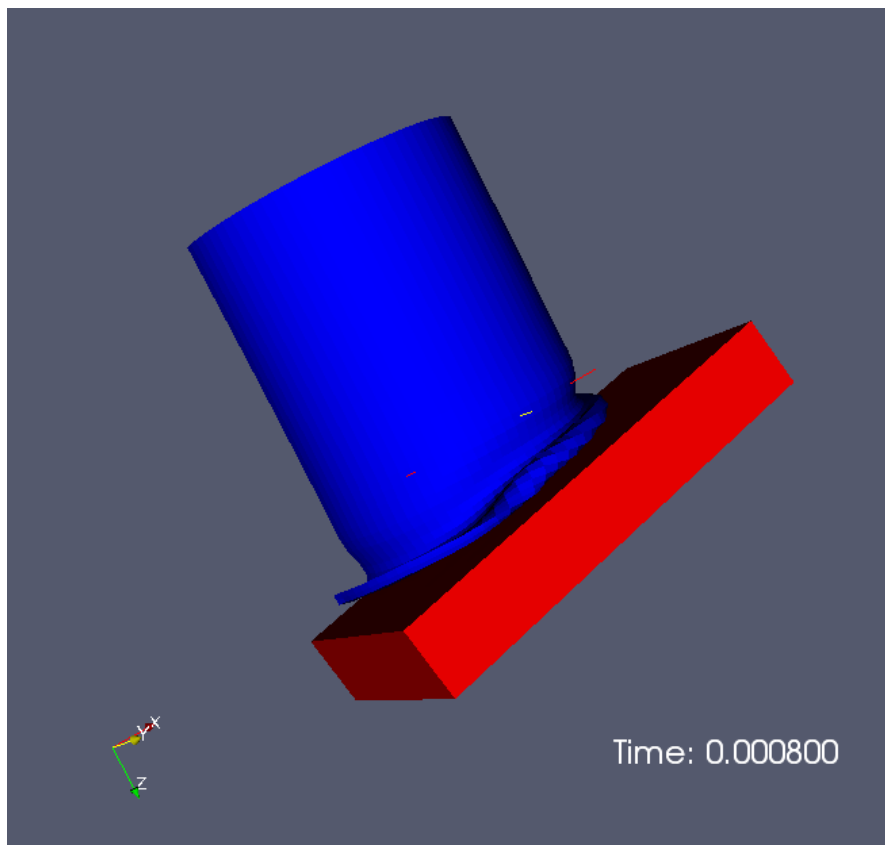


Figure 8. Using the Annotate Time filter to display the current time in a time-varying data set

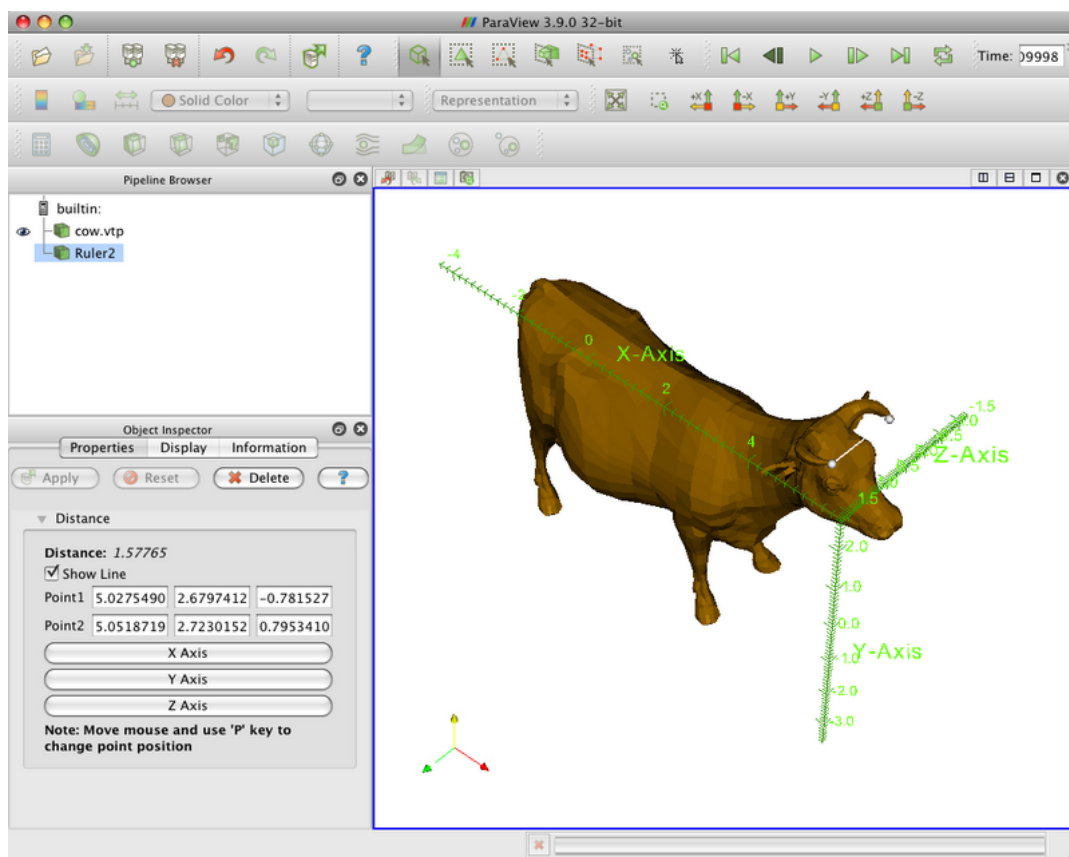
## Other sources as Annotations

Most of the other items in the **Sources** menu can also be used for annotation of the 3D scene. Placement of these is similar to placement of the 3D Text source. In particular the **2D Glyph**, **Ruler** and **Outline** sources are frequently used for annotation.

The ruler as the name applies allows for some quantitative measurement of the data. The ruler produces a Line widget in the 3D scene. It is most effective to use the 'P' key to place the two ends of the line in the 3D scene in turn. When placed the Properties tab for the ruler lists the length of the line you've placed in space. In **Figure 9** A ruler is being used to measure the distance between the cows horns.

## Cube Axes

Finally in the **Annotation** area of the **Display** tab of any object shown in a 3D view you can toggle the display of a **Cube Axes Display**. Like the ruler source, the cube axes allows you to inspect to world space extent of objects in the scene. When this display is enabled, a gridded, labelled bounding box is drawn around the object that lets you quickly determine the world space extent in each dimension of the object.



**Figure 9. Cube Axes and Rulers used to measure a cow and its horns.**

Pressing the **Edit** button next to the enable checkbox brings up a dialog that lets you customize the presentation of the axes display. You have one tab on this dialog for each of the three axes. Use the **Title** entry to specify a label. Use the **Show Axes** checkbox to enable or disable the given axes. **Show Ticks** and **Show Minor Ticks** let you customize the division markers along the axis. **Show Grid Lines** extends the major tick marks so that they completely surround the object. The **Custom Bounds** entry allows you to specify your own extent for that axis label.

At the bottom of the dialog box are controls over all three axes. The **Fly Mode** dropdown controls the way the axes are drawn on different axes dependent on the camera orientation. **Outer Edges** draws opposing edges so as to space the annotation as widely as possible. **Closest Triad** and **Furthest Triad** draw the nearest and furthest corners respectively. The two static options lock the choice of edges to draw so that they do not change as the camera

moves. **Tick Location** draws the tick marks inward toward the center of the object, outward away from the center, or in both directions simultaneously. **Corner Offset** puts some space between the actual bounds and the displayed axes, which is again useful for minimizing visual clutter. Finally you have control over the color of the axes through the **Set Axes Color...** popup color selector.

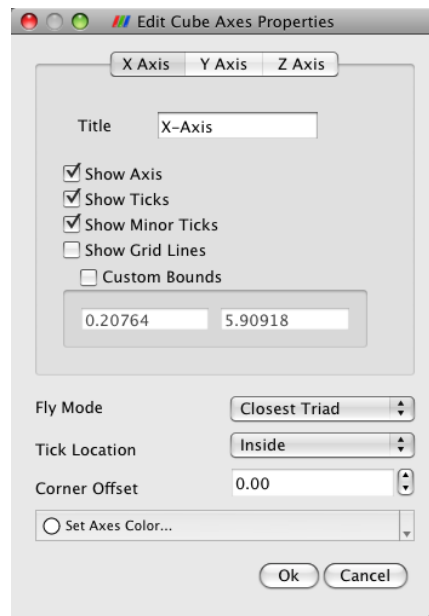


Figure 10. Edit Cube Axes Dialog

## References

- [1] [http://paraview.org/Wiki/ParaView/Displaying\\_Data](http://paraview.org/Wiki/ParaView/Displaying_Data)

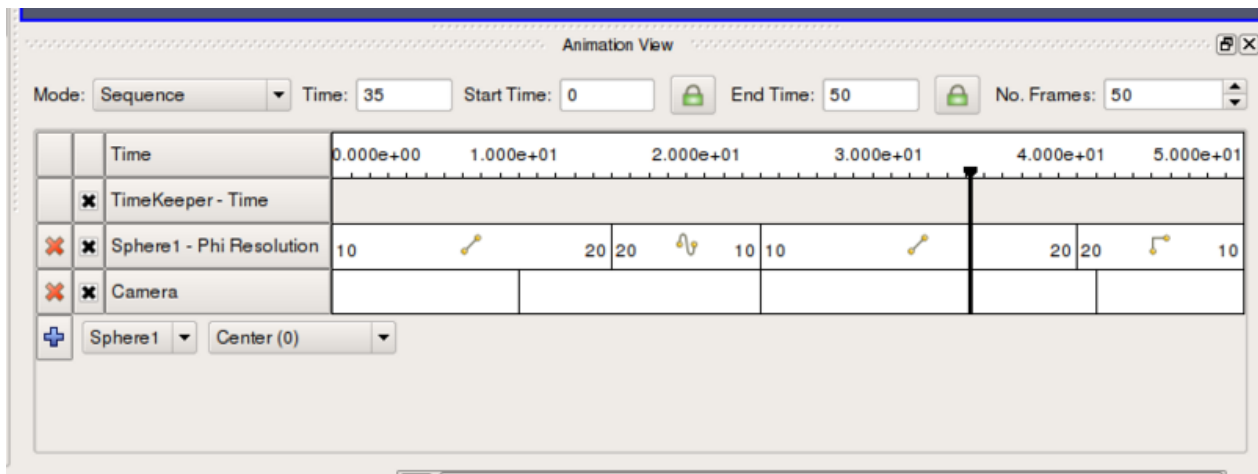
# Animation

## Animation View

In ParaView, you can create animations by recording a series of keyframes. At each keyframe you set values for the properties of the readers, sources and filters that make up the visualization pipeline, as well as the position and orientation of the camera. Once you have chosen the parameters of the animation, you can play through the animation. When you play the animation, you can cache the geometric output of the visualization pipeline in memory. When you subsequently replay the animation, playback will be much faster, because very little computation must be done to generate the images. Also the results of the animation can be saved to image files (one image per animation frame) or to a movie file. The geometry rendered at each frame can also be saved in ParaView's PVD file format, which can be loaded back into ParaView as a time varying data set.

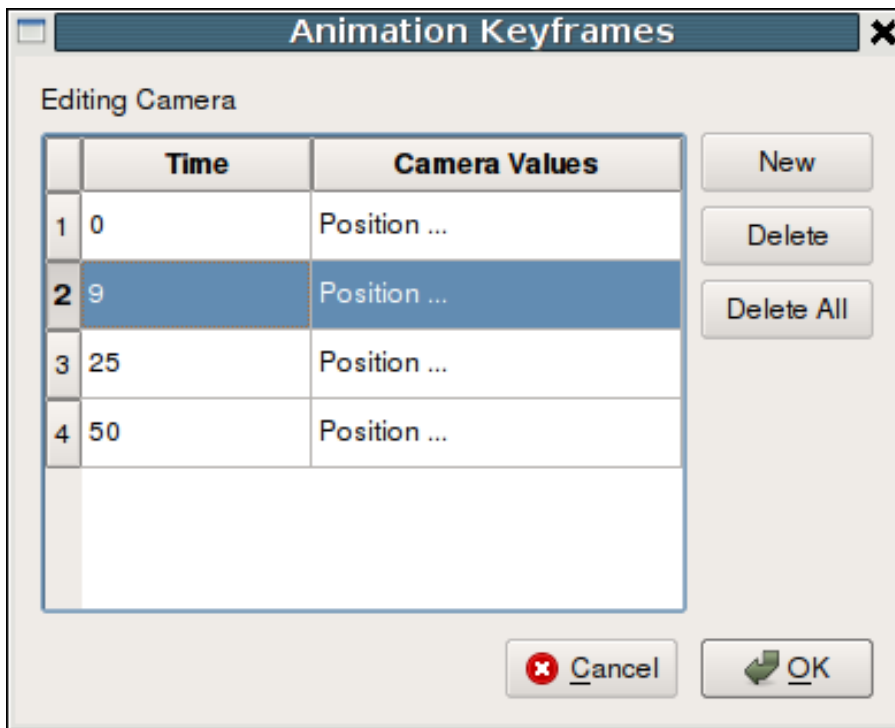
## Animation View

**Animation View** is the user-interface used to create animations by adding keyframes. It is modeled similar to popular animation and key-frame editing application with ability to create tracks for animating multiple parameters. The **Animation View** is accessible from the **View** menu.



This view is presented as a table. Above the table are the controls that control how time progresses in the animation. These were discussed previously. Within the table the tracks of the animation appear as rows, and animation time is presented increasing from left to right. The first row in the table, simply labeled Time, shows the total span of time that the animation can cover. The currently displayed time is indicated both in the Time field at the top and with a thick, vertical, draggable line within the table.

Along the left side of the Animation View is an expandable list of the names of the animation tracks (i.e., a particular object and property to animate). You choose a data source and then a particular property of the data source in the bottom row. To create a animation track with keyframes for that property, click the plus mark on the left-hand side. This will create a new track. In the figure, tracks already exist for SphereSource1's Phi Resolution property and for the Camera's position. To delete a track press the 'X' button. You can temporarily disable a track by un-checking the checkbox on the right of the track. To enter values for the property, double-click within the white area to the right of the track name. This will bring up the Animation Keyframes dialog. Double-clicking in the camera entry brings up a dialog like the one in the following figure.

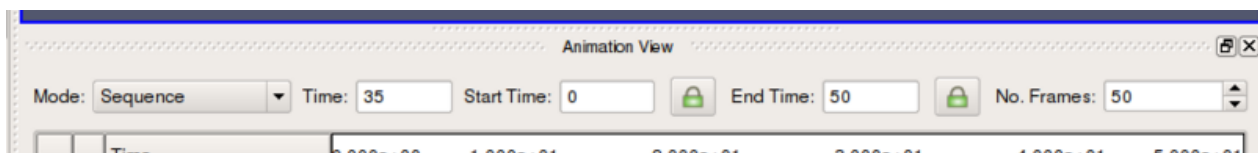


From the Animation Keyframes dialog you can press **New** to create new keyframes or press **Delete** or **Delete All** to delete some or all of them. Clicking **New** will add a new row to the table. In any row, you can click within the Time column to choose a particular time for the keyframe and click in the right-hand column to enter values for the parameter. The exact user interface components that let you set values for the property at the keyframe time vary. When available, you can change the interpolation between two keyframes by double-clicking on the central interpolation column.

Within the tracks of the Animation View, the place in time where each keyframe occurs is shown as a vertical line. The values chosen for the property at that time and the interpolation function used between that value and the next are shown as text when appropriate. In previous figure for example, the sphere resolution begins at 10 and then changes to 20 varying by linear interpolation between them. The camera values are too lengthy to show as text so they are not displayed in the track, but we can easily see that there are four keyframes spaced throughout the animation. The vertical lines in the tracks themselves are draggable, so you can easily adjust the time at which each keyframe occurs.

## Animation View Header

The **Animation View** has a header-bar that lets you control some properties of the animation itself.



**Mode** controls the animation playback mode. ParaView supports 3 modes for playing animation. In **Sequence** mode, the animation is played as a sequence of images (or frames) generated one after the other and rendered in immediate succession. The number of frames is controlled by the **No. Frames** spinbox at the end of the header. Note that the frames are rendered as fast as possible. Thus the viewing frame rate depends on the time needed to generate and render each frame.

In **Real Time** mode, the **Duration** spinbox (replacing the **No. Frames** spinbox) indicates the time in seconds over which the entire animation should run. Each frame is rendered using the current wall clock time in seconds relative to the start time. The animation runs for nearly the number of seconds specified by the Duration (secs) spinbox. In

turn, the number of frames actually generated (or rendered) depends on the time to generate (or render) each frame.

In **Snap To TimeSteps** mode, the number of frames in the animation is determined by the number of time values in the data set being animated. This is the animation mode used for ParaView's default animations: playing through the time values in a data set one after the other. Default animations are created by ParaView when a data set with time values is loaded; no action is required by the user to create the animation. Note that using this mode when no time-varying data is loaded will result in no animation at all.

In **Sequence** mode, the final item in the header is the **No. Frames** spinbox. This spinbox lets you pick the total number of frames for the animation. Similarly in **Real Time** mode, the final line lets you choose the duration of the animation. In **Snap To Timesteps** mode, the total number of frames is dictated by the data set, and therefore the spinbox is disabled.

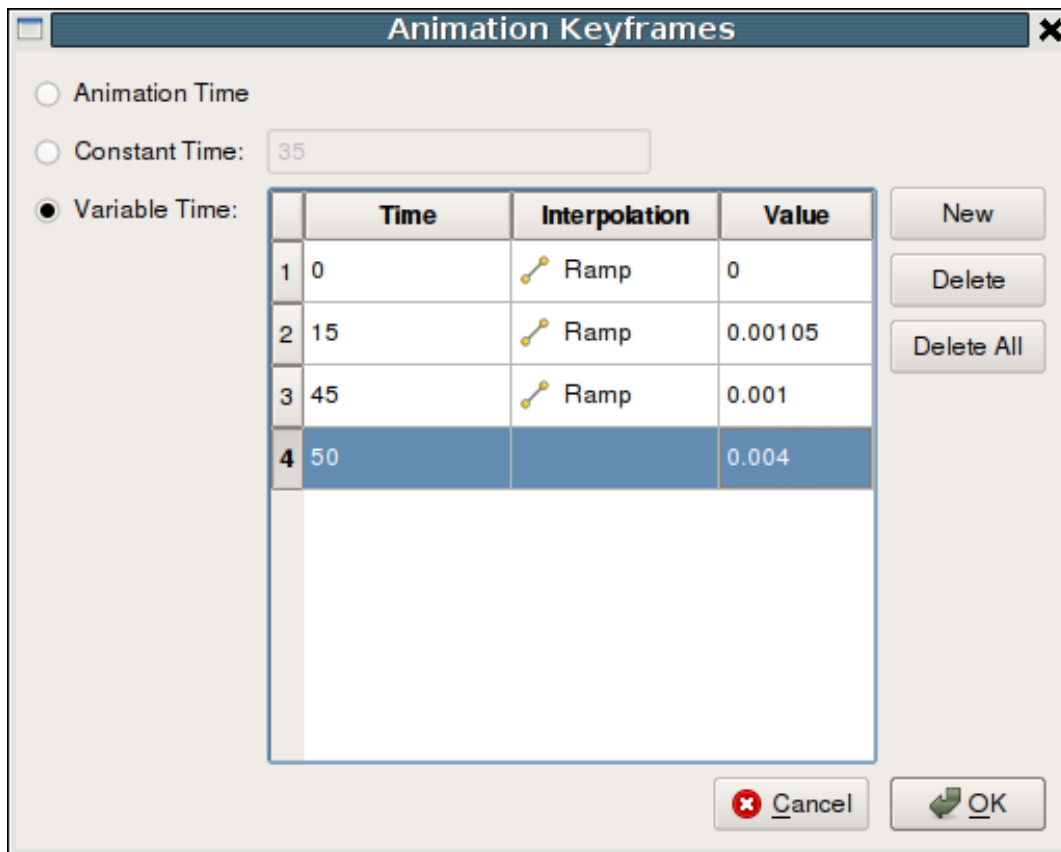
The **Time** entry-box shows the current animation time which is same as shown by a vertical marker in this view. One can change the current animation time by either entering a value in this box, if available or dragging the vertical marker. The **Start Time** and **End Time** entry-boxes display the start and end times for the animation. By default, when you load time varying data sets, the start and end times are automatically adjusted to cover the entire time range present in the data. The lock check-buttons to the right of the **Start Time** and **End Time** widgets will prevent this from happening so that you can ensure that your animation covers a particular time domain of your choosing.

## Animating Time-Varying Data

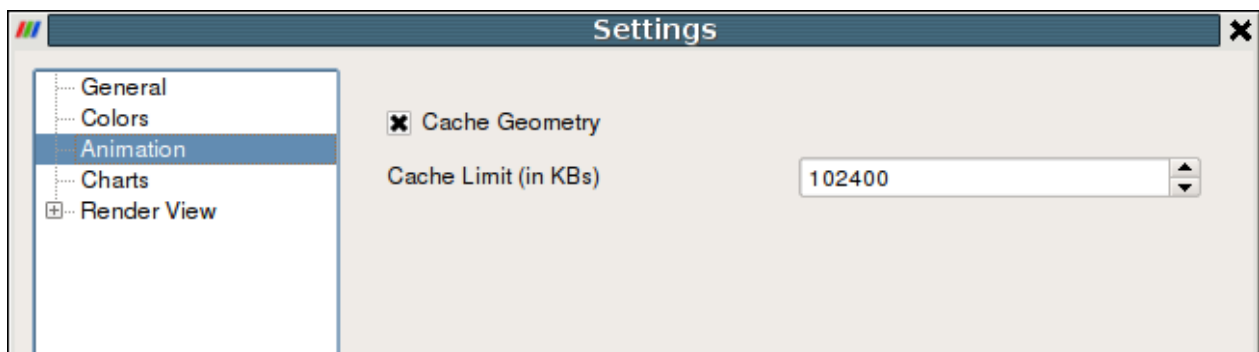
When you load time-varying data, ParaView automatically creates a default animation that allows you to play through the temporal domain of the data without manually creating an animation to do so. With the **Animation View**, you can uncouple the data time from the animation time so that you can create keyframes that manipulate the data time during animation as well.

If you double click in the **TimeKeeper – Time** track, the **Animation Keyframes** dialog, an example of which is shown in Figure below, appears. In this dialog, you can make data time progress in three fundamentally different ways. If the **Animation Time** radio-button is selected, then the data time will be tied to and scaled with the animation time so that as the animation progresses you will see the data evolve naturally. You can select **Constant Time** instead if you want to ignore the time varying nature of the data. In this case you choose a particular time value at which the data will be displayed for the duration of the animation. Finally, you can select the **Variable Time** radio-button to have full control over data time and control it as you do any other animatable property in the visualization pipeline. In the example shown in Figure below, time is made to progress forward for the first 15 frames of the animation, then backward for the next 30, and finally forward for the final 15.





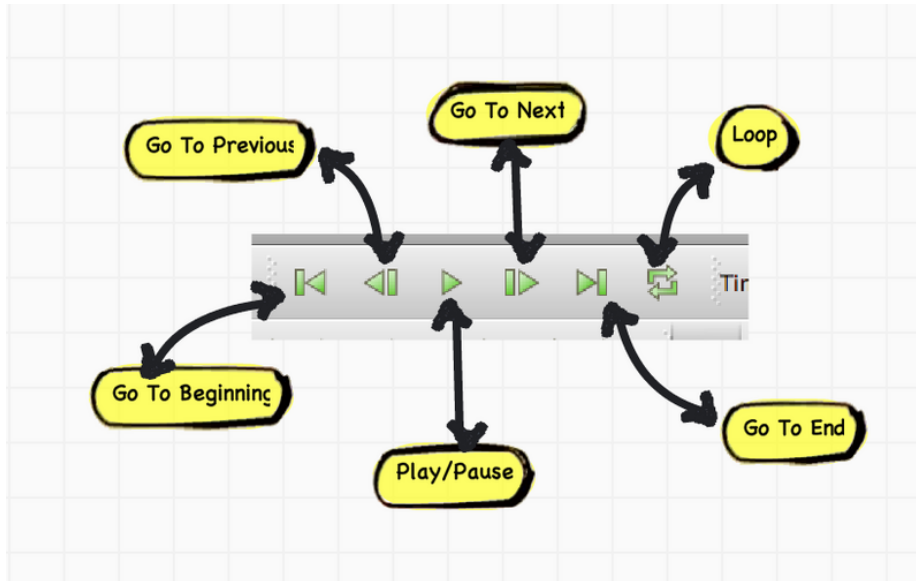
## Animation Settings



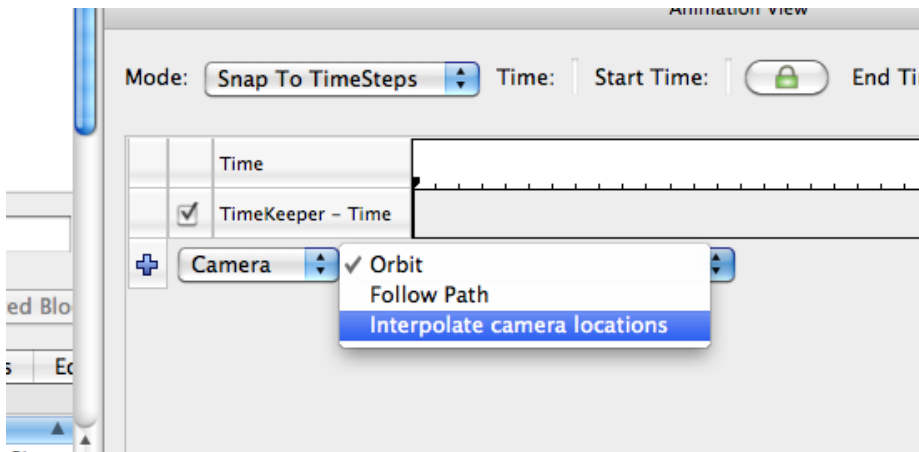
Additional animation properties can be changed using the **Animation** page in the **Edit | Settings** dialog. Using these settings one can control whether geometry must be cached to improve playback performance during looping as well as the maximum size of geometry to cache to avoid memory overflows.

## Playing an Animation

Once you have designed your animation, you can play through it with the VCR controls toolbar.



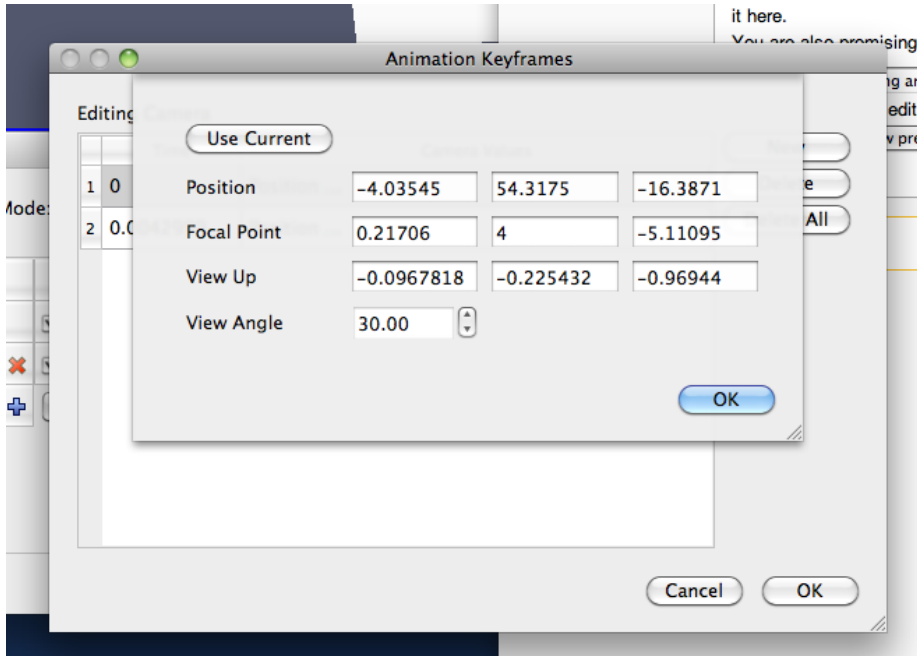
## Animating the Camera



Just like one can change parameters on sources and filters in an animation, one can also change the camera parameters. One can add animation tracks to animate the camera for all the 3D render views in the setup separately. To add a camera animation track for a view, with the view selected, click on the "+" button after choosing "Camera" from the first drop down. The second drop down allows users to choose how to animate the camera. There are three possible options each of which provides different mechanisms to specify the keyframes. It's not possible to change the mode after the animation track has been added, but one can simply delete the track and create a new one.

## Interpolate Camera Locations

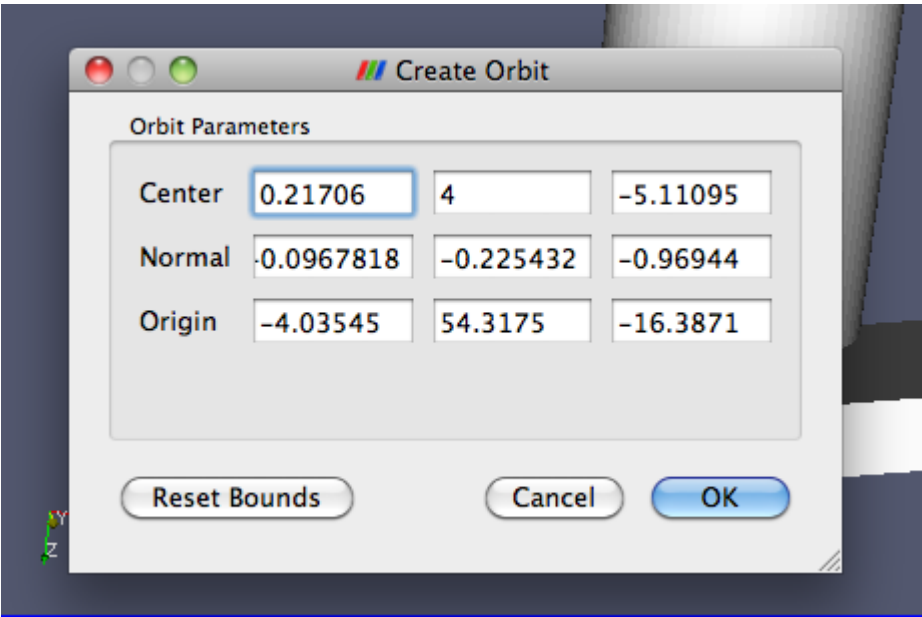
In this mode, the user specifies camera position, focal point, view angle and up direction at each keyframe. The animation player interpolates between these specified locations. As with other parameters, to edit the keyframes, one double-clicks on the track. It is also possible to capture the current location as a keyframe by using the "Use Current" button.



It can be quite challenging to add key frames correctly and frequently to ensure that the animation results in a smooth visualization using this mode.

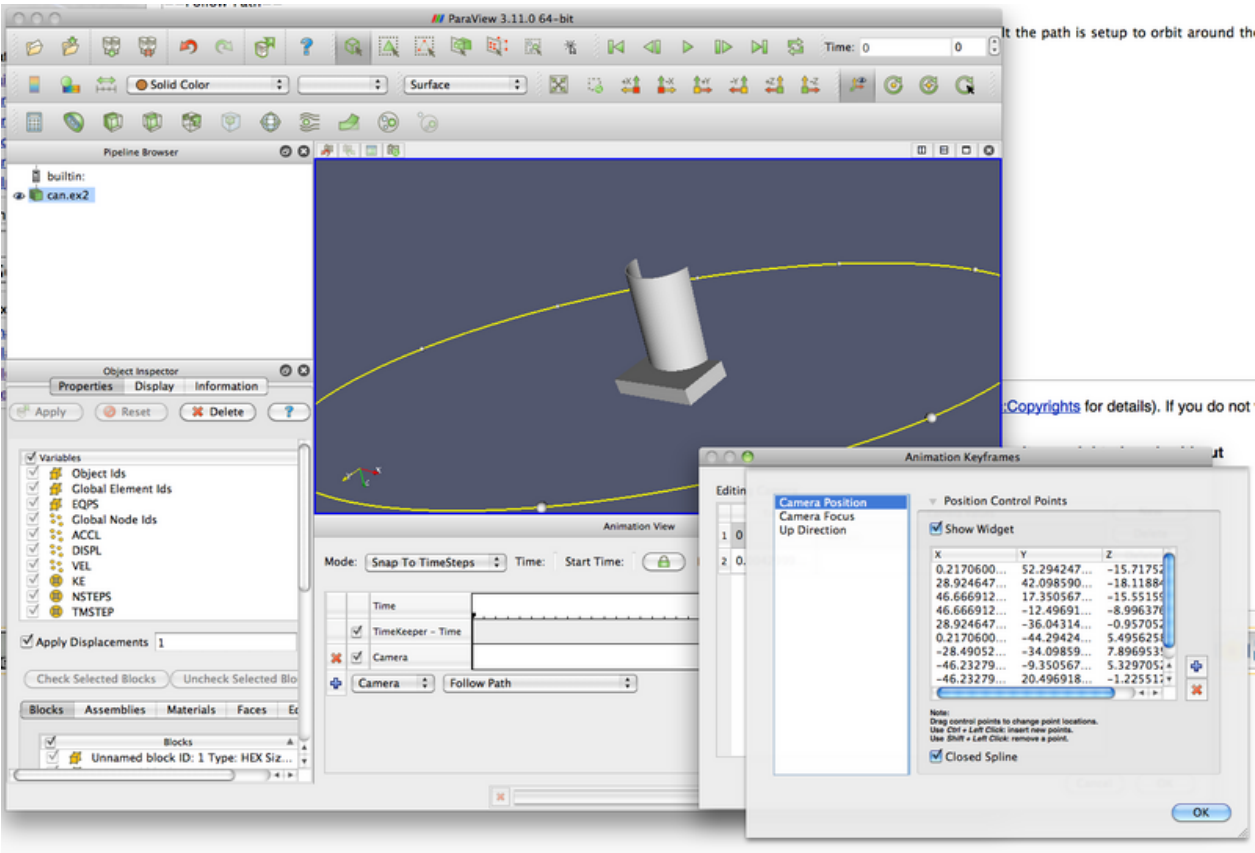
## Orbit

This mode makes it possible to quickly create a camera animation in which the camera revolves around an object or objects of interest. Before adding the Camera track, select the object or objects in the pipeline browser that you want to revolve around. Then choose "Orbit" from the "Camera" combo-box in the Animation View and hit "+". This will popup a dialog where one can edit the orbit parameters such as the center of revolution, normal for the plane of revolution and the origin i.e. a point on the plane of revolution where the revolution begins. By default, the Center is the center of the bounds of the selected objects, Normal is the current up direction used by the camera while the origin is the current camera position.



Follow Path

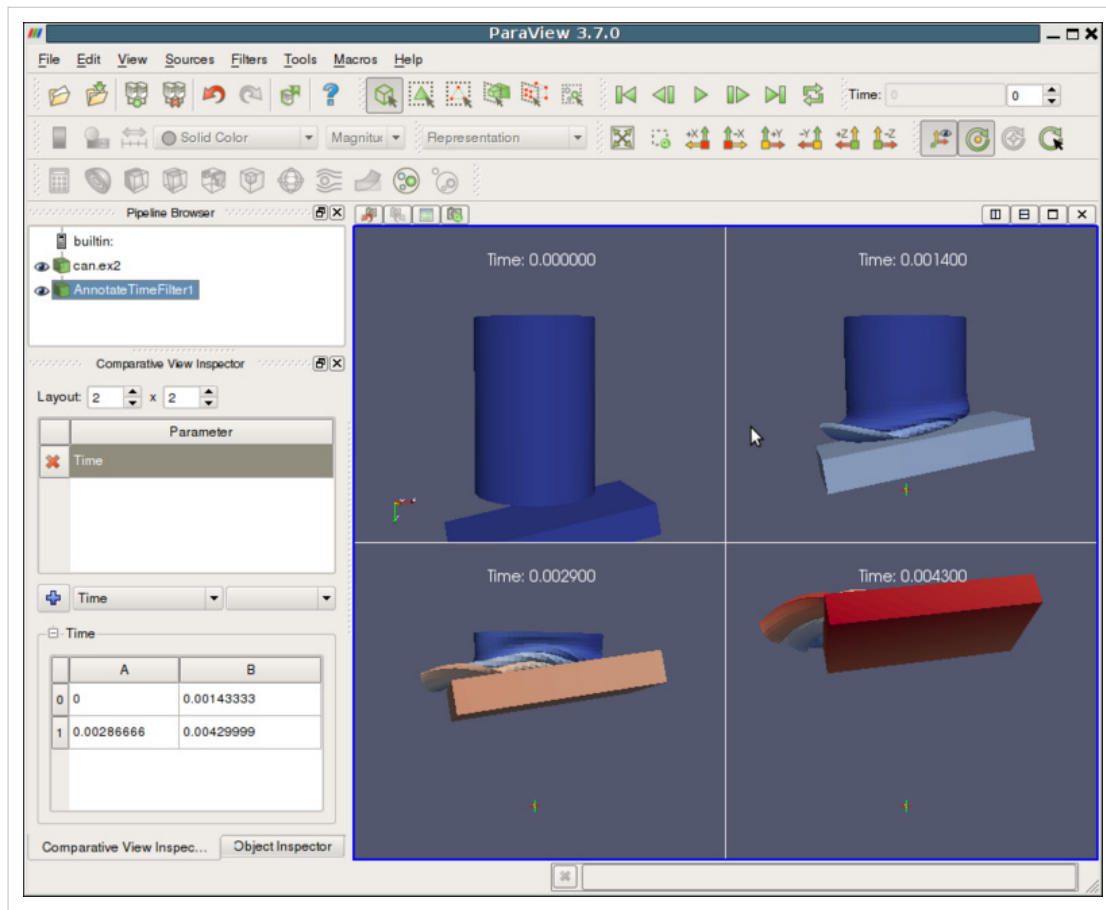
In this mode the user gets the opportunity to specify the path taken by the camera position and camera focal point. By default the path is setup to orbit around the selected objects. User can then edit the keyframe to change the paths. Figure shows the dialog for editing these paths for a keyframe. When "Camera Position" or "Camera Focus" is selected, a widget is shown in the 3D view that can be used to set the path. Use *Ctrl+Left Click* to insert new control points, and *Shift+Left Click* to remove control points. One can also toggle when path should be closed or not.



# Comparative Visualization

## Comparative Views

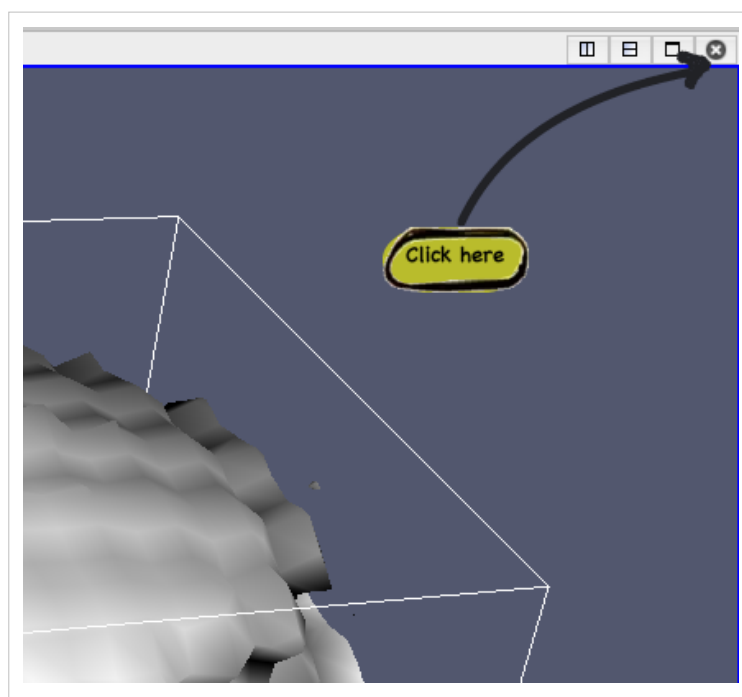
### Introduction



ParaView provides a collection of views that can be used to display a set of visualization on a regular grid. These views are referred to as "Comparative" Views. Comparative Views can be used to display results from a parameter study or to perform parameter studies within ParaView.

## Quick Start

We will start with a short tutorial. Later sections will describe Comparative Views in more detail. First start a Wavelet source. Next apply a Contour filter. Then close the default 3D view by clicking on the small x on the upper right corner of the view.



This should bring up an empty view with several buttons for selecting the view type. Select 3D View (Comparative). Now you should see a 2x2 grid in the view and it should be empty. Next turn on the visibility of the Contour filter. You should see the output of the Contour filter in all 4 views. Next, we are going to vary the contour value across the grid. Bring up the Comparative View Inspector (View -> Comparative View Inspector). This inspector allows you to configure the parameters of the Comparative View.

**1. Determines the layout for the parameter study grid.**

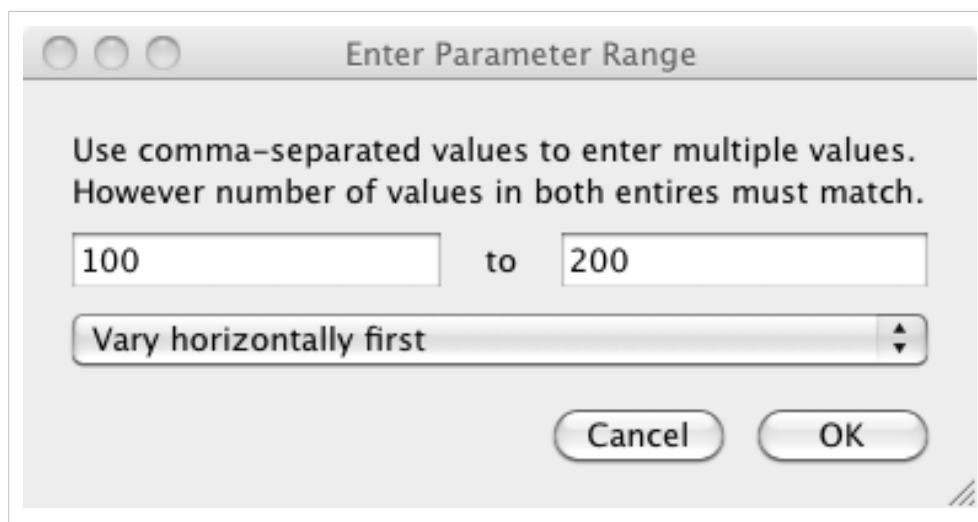
**2. Parameters being studied:**  
Select a parameter here to see its values in the

**3. Select and add new parameters to the study.**

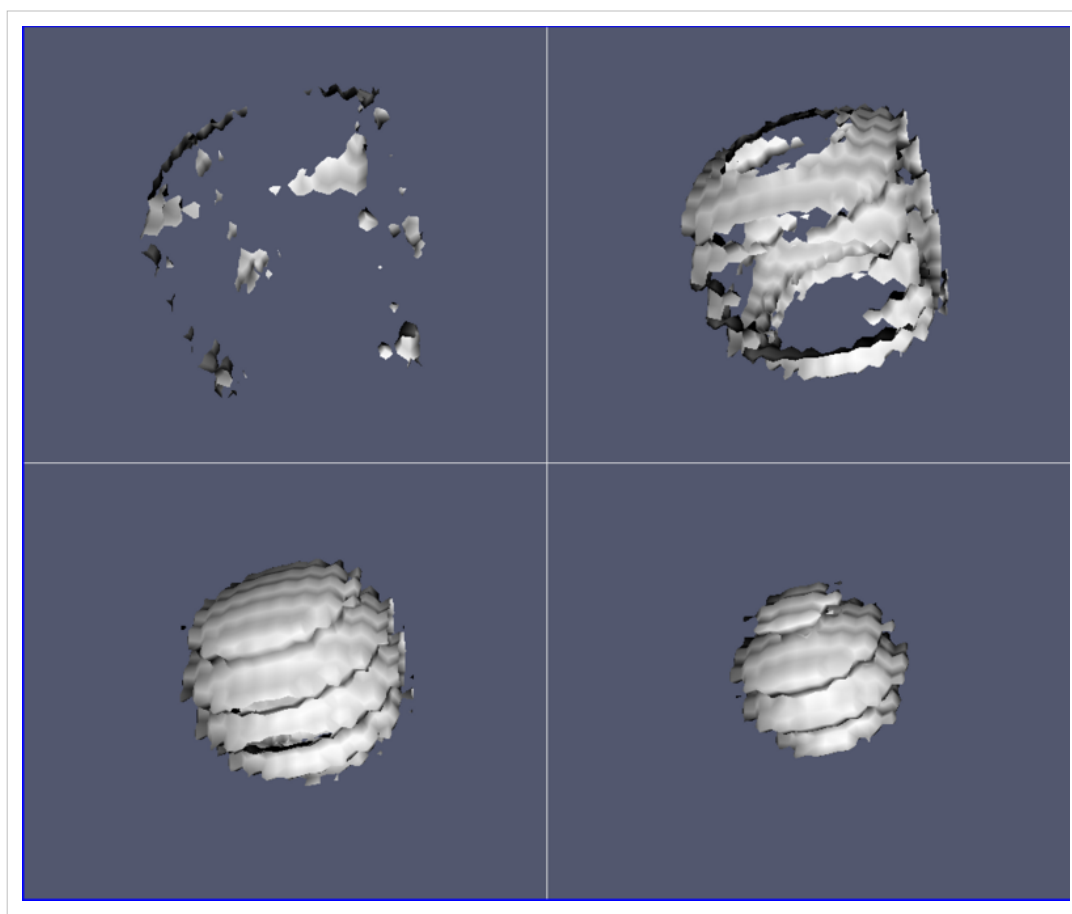
**4. Parameter values over the study grid:**  
 \* Double click on a value to change it.  
 \* Drag and select a range to change the values for those cells.  
 \* When multiple values are accepted by the parameter e.g. Isovalues for the Contour filter, you can enter comma-separated values

| Comparative View Inspector          |            |            |
|-------------------------------------|------------|------------|
| Layout: 2 x 2                       |            |            |
| Parameter                           |            |            |
| <input checked="" type="checkbox"/> | Time       |            |
| <input type="checkbox"/>            |            |            |
| <input type="checkbox"/>            |            |            |
| <input type="checkbox"/>            |            |            |
| <input type="checkbox"/>            | Time       |            |
|                                     | A          | B          |
| 0                                   | 0          | 0.00143333 |
| 1                                   | 0.00286666 | 0.00429999 |

From the drop-down menu select Contour1. The parameter should be set to Isosurfaces. If it is not, change it. Next, in the grid that has the numbers, click on the upper left value and drag to the lower right value. This should bring up the following dialog. Enter 100 and 200 for the two values.



When you click OK, the Comparative View should update and you should see something like the following image.



## View

There are three types of Comparative Views:

- Comparative 3D View
- Comparative Line Chart View
- Comparative Bar Chart View

All of these views contain sub-views laid-out in an m-by-n grid where m and n are determined by the Comparative View Inspector settings. The type of the sub-view depends on the type of the Comparative View, 3D View for Comparative View, Line Chart for Line Chart Comparative View etc. Each sub-view display the output from the same pipeline objects, depending on what is set visible in the Pipeline Inspector. The only things that change from view to view are the parameters that are varied through the Comparative View Inspector. Furthermore, the view settings are synchronized between all sub-views. For example, changing the camera in one 3D sub-view will cause the camera on all other sub-views to change.

**Note:** Not all features of the single views are supported in their comparative siblings. For example, it is not possible to perform surface selection on a Comparative 3D View.

## Comparative View Inspector

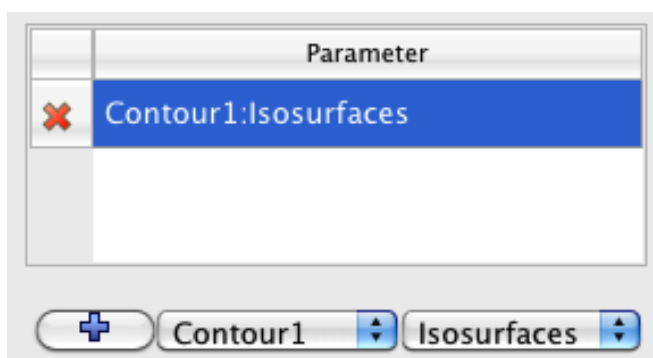
The Comparative View Inspector (View -> Comparative View Inspector) is where you can configure all of the parameters of Comparative Views. Note that if you have more than one Comparative View, the Inspector will change the setting of the active one. Below we describe various parts of the Comparative View Inspector.

### Layout



The layout widget allows you to configure the number of sub-views within a Comparative View. The first value controls how many cells there are in the horizontal direction whereas the second value controls how many cells there are in the vertical direction. Note that if you already setup one or more parameters to vary in the view, the Inspector will try to maintain your values when you adjust the size as much as possible. If you manually entered any individual values, they will not change when you add more rows or columns. On the other hand, all values that have been automatically computed based on a range will be updated as the number of cells change.

### Parameter Selection

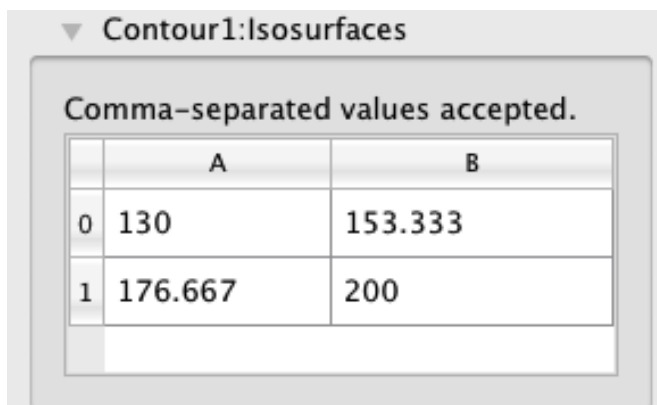


This widget allows you to add new parameters to vary and also to select a property if more than one is available. To add a property, first select the pipeline object from the first menu, then select its parameter from the second value. Once you are done with the selection, click on the + button to add the parameter. This parameter will now show in the list of parameters. To delete a parameter from the list, click on the corresponding x on the left side of the parameter selection widget. Note that once you add a new parameter, ParaView will try to assign good default values



to each cell. For example, if you add Contour:Isosurfaces, ParaView will assign values ranging from minimum scalar value to maximum scalar value.

## Editing Parameter Values



You can edit the parameter values using the spreadsheet widget. You can either:

- Change individual value by double clicking on a cell and editing it or
- Change a group of cells by clicking on the first one and dragging to the last one. ParaView will then ask you to enter minimum and maximum values. If you selected cells that span more than one direction, it will also ask you to choose which way values vary.

Let's examine the range selection a bit further. Say that you selected a 2x2 area of cells and entered 0 for the minimum and 10 for the maximum. If you select **Vary Horizontally First**, the values will be:

|      |      |
|------|------|
| 0    | 3.33 |
| 6.66 | 10   |

If you select, **Vary Vertically First**, the values will be

|      |      |
|------|------|
| 0    | 6.66 |
| 3.33 | 10   |

If you select, **Vary Only Horizontally**, the value will be

|   |    |
|---|----|
| 0 | 10 |
| 0 | 10 |

If you select, **Vary Only Vertically**, the value will be

|    |    |
|----|----|
| 0  | 0  |
| 10 | 10 |

The last two options may sound useless. Why have multiple cells with the same values? However, if you consider that more than one parameter can be varied in a Comparative View, you will realize how useful they are. For example, you can change parameter A horizontally while varying parameter B vertically to create a traditional spreadsheet of views.

## **Performance**

### **Computational**

ParaView will run the pipeline connected to all visible pipeline objects for each cell serially. Therefore, the time to create a Comparative Visualization of N cells should be on the order of N times the time to create the visualization of 1 cell.

### **Memory**

ParaView will only store what is needed to display the results for each cell except the last one. The last cell will contain the representation as well as the full dataset, same as any single view. For example, when using the surface representation, the total memory used will be the total of memory used by the geometry in each cell plus the memory used by the dataset of the last cell.

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# Remote and Parallel Large Data Visualization

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## Parallel ParaView

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### Parallel ParaView

One of the main purposes of ParaView is to allow users to create visualizations of large data sets that reside on parallel systems without first collecting the data to a single machine. Transferring the data is often slow and wasteful of disk resources, and the visualization of large data sets can easily overwhelm the processing and especially memory resources of even high-performance workstations. This chapter first describes the concepts behind the parallelism in ParaView. We then discuss in detail the process of starting up ParaView's parallel server components. Lastly we explain explains how a parallel visualization session is initiated from within the user interface. Parallel rendering is an essential part of parallel ParaView. So essential that we've given it its own chapter in this version of the book.

The task of setting up a cluster for Visualization is unfortunately outside of the scope of this book. However there are several online resources that will help to get you started including:

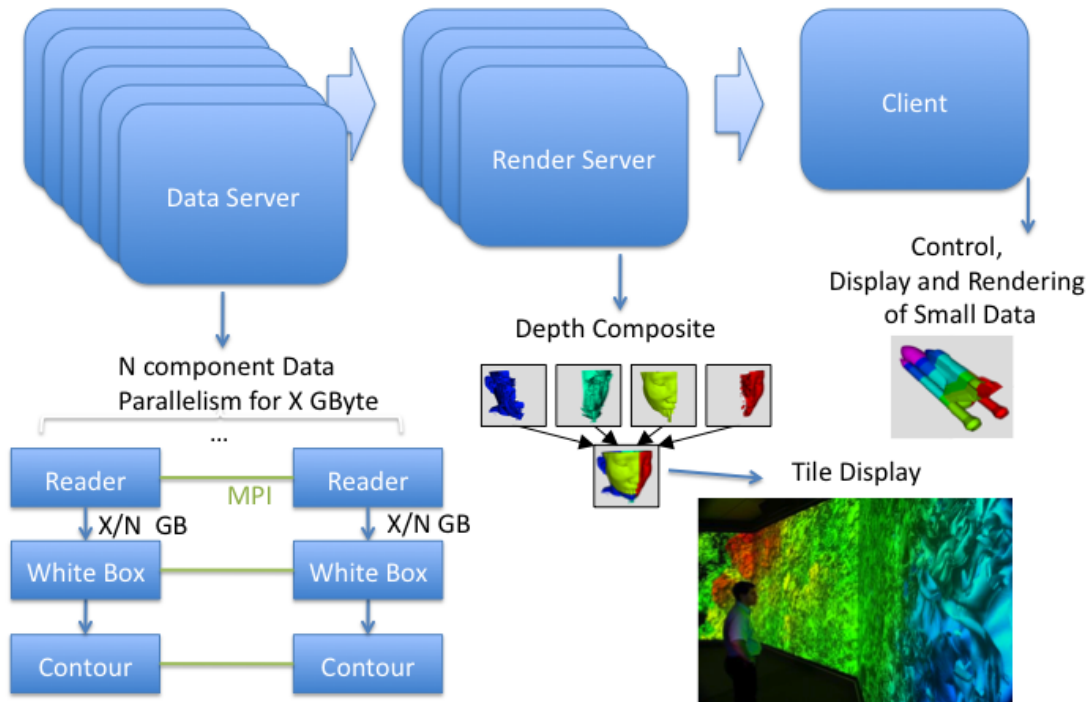
- [http://paraview.org/Wiki/Setting\\_up\\_a\\_ParaView\\_Server](http://paraview.org/Wiki/Setting_up_a_ParaView_Server),
- <http://www.iac.es/sieinvens/siepedia/pmwiki.php?n=HOWTOs.ParaviewInACluster>

and

- [http://paraview.org/Wiki/images/a/a1/Cluster09\\_PV\\_Tut\\_Setup.pdf](http://paraview.org/Wiki/images/a/a1/Cluster09_PV_Tut_Setup.pdf)
-

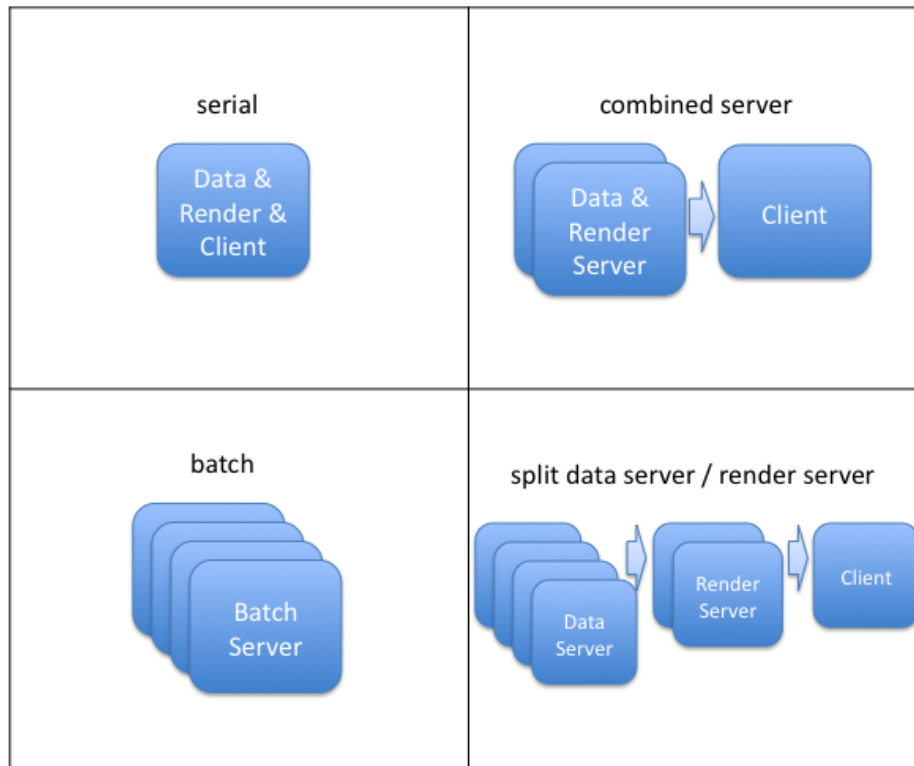
## Parallel Structure

ParaView has three main logical components: **client**, **data server**, and **render server**. The client is responsible for the GUI and is the interface between you the user and ParaView as a whole. The data server reads in files and processes the data through the pipeline. The render server takes the processed data and renders it to present the results to you.



**Figure 1. Parallel Architecture**

The three logical components can be combined in various different configurations. When ParaView is started, the client is connected to what is called the builtin server; in this case all three components exist within the same process. Alternatively, you can run the server as an independent program and connect a remote client to it, or run the server as a standalone parallel batch program without a GUI. In this case the server process contains both the data and render server components. The server can also be started as two separate programs: one for the data server and one for the render server. The server programs are data-parallel programs that can be run as a set of independent processes running on different CPUs. The processes use MPI to coordinate their activities as each works on different pieces of the data.



**Figure 2. Common Configurations of the logical components**

## Client

The client is responsible for the user interface of the application. ParaView's general-purpose client was written to make powerful visualization and analysis capabilities available from an easy-to-use interface. The client component is a serial program that controls the server components through the Server Manager API.

## Data Server

The data server is primarily constructed from VTK readers, sources, and filters. It is responsible for reading and/or generating data, processing it, and producing geometric models that the render server and client will display. The data server exploits data parallelism by partitioning the data, adding ghost levels around the partitions as needed, and running synchronous parallel filters. Each data server process has an identical VTK pipeline, and each process is told which partition of the data it should load and process. By splitting the data ParaView is able to use the entire aggregate system memory and thus make large data processing possible.

## Render Server

The render server is responsible for rendering the geometry. Like the data server, the render server can be run in parallel and has identical visualization pipelines (only the rendering portion of the pipeline) in all of its processes. Having the ability to run the render server separately from the data server allows for an optimal division of labor between computing platforms. Most large computing clusters are primarily used for batch simulations and do not have hardware rendering resources. Since it is not desirable to move large data files to a separate visualization system, the data server can run on the same cluster that ran the original simulation. The render server can be run on a separate visualization cluster that has hardware rendering resources.

It is possible to run the render server with fewer processes than the data server but never more. Visualization clusters typically have fewer nodes than batch simulation clusters, and processed geometry is usually significantly smaller

than the original simulation dump. ParaView repartitions the geometric models on the data server before they are sent to the render server.

## MPI Availability

Until recently, in order to use ParaView's parallel processing features, one needed to build ParaView from the source code as described in the Appendix. This was because there are many different versions of MPI, the library ParaView's servers use internally for parallel communication, and for high performance computer users it is extremely important to use the version that is delivered with your networking hardware.

As of ParaView 3.10 however, we have begun to package MPI with our binary releases. If you have a multi-core workstation, you can now simply turn on the **Use Multi-Core** setting under ParaView's Settings to make use of all of them. This option makes parallel data server mode the default configuration, which can be very effective when you are working on computationally bound intensive processing tasks.

Otherwise, and when you need to run ParaView on an actual distributed memory cluster, you need to start up the various components and establish connections between them as is described in the next section.

# Starting the Server(s)

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## Client / Server Mode

Client / Server Mode refers to a parallel ParaView session in which data server and render server components reside within the same set of processes, and the client is completely separate. The `pvserver` executable combines the two server components into one process.

You can run `pvserver` as a serial process on a single machine. If ParaView was compiled with parallel support, you can also run it as an MPI parallel program on a group of machines. Instructions for starting a program with MPI are implementation- and system-dependent, so contact your system administrator for information about starting an application with MPI. With the MPICH implementation of MPI, the command to start the server in parallel usually follows the format shown here.

```
mpirun -np number_of_processes path_to/pvserver arguments for pvserver
```

By default, `pvserver` will start and then wait for the client to connect to it. See the next section for a full description. Briefly, to make the connection, select **Connection** from the **File** menu, select (or make and then select) a configuration for the server, and click **Connect**. Note that you must start the server before the client attempts to connect to it.

If the computer running the server is behind a firewall, it is useful to have the server initiate the connection instead of the client. The **--reverse-connection** (or **-rc**) command-line option tells the server to do this. The server must know the name of the machine to which it should connect; this is specified with the **--client-host** (or **-ch**) argument. Note that when the connection is reversed, you must start the client and instruct it to wait for a connection before the server attempts to connect to it.

The client-to-server connection is made over TCP, using a default port of 11111. If your firewall puts restrictions on TCP ports, you may want to choose a different port number. In the client dialog, simply choose a port number in the **Port** entry of the **Configure New Server** dialog. Meanwhile, give `pvserver` the same port number by including **--server-port** (or **-sp**) in its command-line argument list.

An example command line to start the server and have it initiate the connection to a particular client on a particular port number is given below.

```
pvserver -rc -ch=magrathea -sp=26623
```

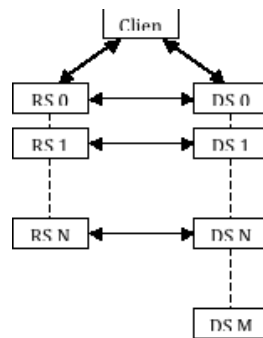
---

## Render/Server Mode

The render server allows you to designate a separate group of machines (i.e., apart from the data server and the client) to perform rendering. This parallel mode lets you use dedicated rendering machines for parallel rendering rather than relying on the data server machines, which may have limited or no rendering capabilities. In ParaView, the number of machines (N) composing the render server must be no more than the number (M) composing the data server.

Note that although it is true that at some large installations with particularly high performing parallel rendering resources it can be very efficient to run ParaView in Render/Server Mode, we have found in practice that it is almost always the case that the data transfer time between the two servers overwhelms the speedup gained by rendering on the dedicated graphics cluster. For this reason we typically recommend that you combine the data and render servers together as one component and either render in software via Mesa on the data processing cluster or do all of the visualization processing directly on the GPU cluster.

If you still want to break up the data processing and rendering tasks, there are two sets of connections that must be made for ParaView to run in render-server mode. The first connection set is between the client and the first node of each of the data and render servers. The second connection set is between the nodes of the render server and the first N nodes of the data server. Once all of these connections are established, they are bi-directional. The diagram in Figure 5 depicts the connections established when ParaView is running in render server mode. Each double-ended arrow indicates a bi-directional TCP connection between pairs of machines. The dashed lines represent MPI connections between all machines within a server. In all the diagrams in this section, the render server nodes are denoted by RS 0, RS 1, ..., RS N. The data server nodes are similarly denoted by DS 0, DS 1, ..., DS N, ..., DS M.



**Figure 5. Connections required in render server mode**

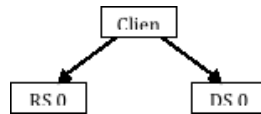
The establishment of connections between client and servers can either be forward (from client to servers) or reverse (from servers back to client). Likewise, the connections between render server and data server nodes can be established either from the data server to the render server or from the render server to the data server.

The main reason for reversing the direction of any of the initial connections is that machines behind firewalls are able to initiate connections to machines outside the firewall, but not vice versa. If the data server is behind a firewall, the data server should initiate the connection with the client, and the data server nodes should connect to the render server nodes. If the render server is behind the firewall instead, both servers should initiate connections to the client, but now the render server nodes should initiate the connections with the nodes of the data server.

In the remaining diagrams in this section, each arrow indicates the direction in which the connection is initially established. Double-ended arrows indicate bi-directional connections that have already been established. In the example command lines, optional arguments are enclosed in []'s. The rest of this section will be devoted to discussing the two connections required for running ParaView in render server mode.

## Connection 1: Connecting the client and servers

The first connection that must be established is between the client and the first node of both the data and render servers. By default, the client initiates the connection to each server, as shown in Figure 6. In this case, both the data server and the render server must be running before the client attempts to connect to them.

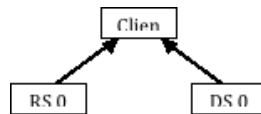


**Figure 6. Starting ParaView in render-server mode using standard connections**

To establish the connections shown above, do the following. First, from the command line of the machine that will run the data server, enter “pvdataserver” to start it. Next, from the command line of the machine that will run the render server, enter “pvrenderserver” to start the render server. Now, from the machine that will run the client, start the client application, and connect to the running servers, as described in section 1.2 and summarized below.

Start ParaView and select **Connect** from the **File** menu to open the **Choose Server** dialog. Select **Add Server** to open the **Configure New Server** dialog. Create a new server connection with a **Server Type** of **Client / Data Server / Render Server**. Enter the machine names or IP addresses of the server machines in the **Host** entries. Select **Configure**, and then in the **Configure Server** dialog, choose **Manual**. Save the server configuration, and **Connect** to it. At this point ParaView will establish the two connections. This is similar to running ParaView in client/server mode but with the addition of a render server.

The connection between the client and the servers can also be initiated by the servers. As explained above, this is useful when the servers are running on machines behind a firewall. In this case, the client must be waiting for both servers when they start. The diagram indicating the initial connections is shown in Figure 7.



**Figure 7. Reversing the connections between the servers and the client.**

To establish the connections shown above, start by opening the **Configure New Server** dialog on the client. Choose **Client / Data Server / Render Server (reverse connection)** for the **Server Type** in the **Configure New Server** dialog. Next, add both `--reverse-connection` (or `-rc`) and `--client-host` (or `-ch`) to the command lines for the data server and render server. The value of the `--client-host` parameter is the name or IP address of the machine running the client. You can use the default port numbers for these connections, or you can specify ports in the client dialog by adding the `--data-server-port` (or `-dsp`) and `--render-server-port` (or `-rsp`) command-line arguments to the data server and render server command lines, respectively. The port numbers for each server must agree with the corresponding **Port** entries in the dialog, and they must be different from each other.

For the remainder of this chapter, `-rc` will be used instead of `--reverse-connection` when the connection between the client and the servers is to be reversed.

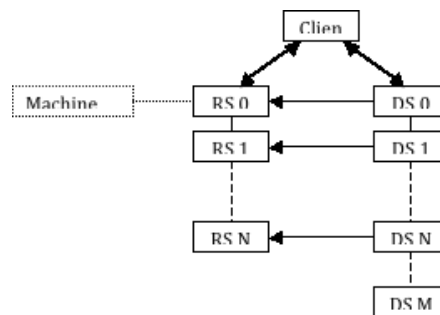


## Connection 2: Connecting the render and data servers

After the connections are made between the client and the two servers, the servers will establish connections with each other. In parallel runs, this server-to-server connection is a set of connections between all N nodes of the render server and the first N nodes of the data server. By default the data server initiates the connection to the render server, but this can be changed with a configuration file. The format of this file is described below.

The server that initiates the connection must know the name of the machine running the other server and the port number it is using. In parallel runs, each node of the connecting server must know the name of the machine for the corresponding process in the other server to which it should connect. The port numbers are randomly assigned, but they can be assigned in the configuration file as described below.

The default set of connections is illustrated in Figure 8. To establish these connections, you must give the data server the connection information discussed above, which you specify within a configuration file. Use the `--machines` (or `-m`) command line argument to tell the data server the name of the configuration file. In practice, the same file should be given to all three ParaView components. This ensures that the client, the render server, and the data server all agree on the network parameters.



**Figure 8. Initializing the connection from the data server to the render server.**

An example network configuration file, called `machines.pvx` in this case, is given below.

```

<?xml version="1.0" ?>
<pvx>

  <Process Type="client">>
  </Process>

  <Process Type="render-server">
    <Option Name="render-node-port" Value="1357"/>
    <Machine Name="rs_m1"
      Environment="DISPLAY=rs_m1:0"/>
    <Machine Name="rs_m2"
      Environment="DISPLAY=rs_m2:0"/>
  </Process>

  <Process Host="data-server">
    <Machine Name="ds_m1" />
    <Machine Name="ds_m2" />
    <Machine Name="ds_m3" />
    <Machine Name="ds_m4" />
  </Process>

</pvx>

```

Sample command-line arguments that use the configuration file above to initiate the network illustrated in Figure 8 are given below.

```
mpirun -np 2 pvdataserver -m=machines.pvx
```

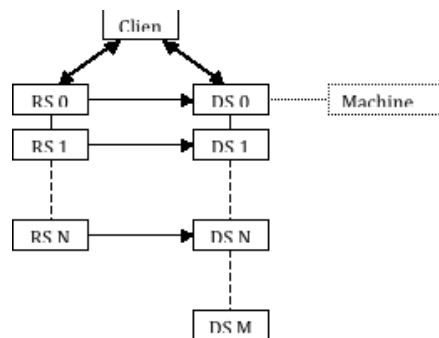
```
mpirun -np 2 pvrenderserver -m=machines.pvx
```

```
paraview -m=machines.pvx
```

It should be noted that the machine configuration file discussed here is a distinct entity from, and has a different syntax from, the server configuration file discussed at the end of the next section. That file is read only by the client; the file discussed here will be given to the client and both servers.

In the machines file above, the render-node-port entry in the render server's XML element tells the render server the port number on which it should listen for a connection from the data server, and it tells the data server what port number it should attempt to contact. This entry is optional, and if it does not appear in the file the port number will be chosen automatically. Note that it is not possible to assign port numbers to individual machines within the server; all will be given the same port number or use the automatically chosen one. Note also that each render server machine is given a display environment variable in this file. This is not required to establish the connections, but it is helpful if you need to assign particular X11 display names to the various render server nodes.

The initial connection between the nodes of the two servers is made from the data server to the render server. You can reverse this such that the render server nodes connect to the corresponding nodes of the data server instead as shown in Figure 9.



**Figure 9. Reversing the connection between the servers and client and connecting the render server to the data server.**

Typically when the server connection is reversed, the direction of the connection between the client and the servers is also reversed (e.g., if the render server is behind a firewall). In this case, the render server must have the machine names and a connection port number to connect to the data server. The same XML file is used for this arrangement as with the standard connection. The only difference in this case is that the render-node-port entry, if it is used, must appear in the data server's XML element instead of the render server's element. Example command-line arguments to initiate this type of network are given here.

```
paraview -m=machines.pvx
```

```
mpirun -np M pvdataserver -m=machines.pvx -rc -ch=client
```

```
mpirun -np N pvrenderserver -m=machines.pvx -rc -ch=client
```

# Connecting to the Server

## Connecting the Client

You establish connections between the independent programs that make up parallel ParaView from within the client's user interface. The user interface even allows you to spawn the external programs and then automatically connect to them. Once you specify the information that ParaView needs to connect to, or spawn and connect to, the server components, ParaView saves it to make it easy to reuse the same server configuration at a later time. Some visualization centers in fact provide predefined configurations which makes it trivial to connect to a server that is tailored to that center by simply choosing one from a list in ParaView's GUI.

The **Choose Server** dialog shown below is the starting point for making and using server configurations. The **Connect** entry on the ParaView client's **File** menu brings it up. The dialog shows the servers that you have previously configured. To connect to a server, click its name to select it; then click the **Connect** button at the bottom of the dialog box. To make changes to settings for a server, select it, and click **Edit Server**. To remove a server from the list, select it, and click **Delete Server**.

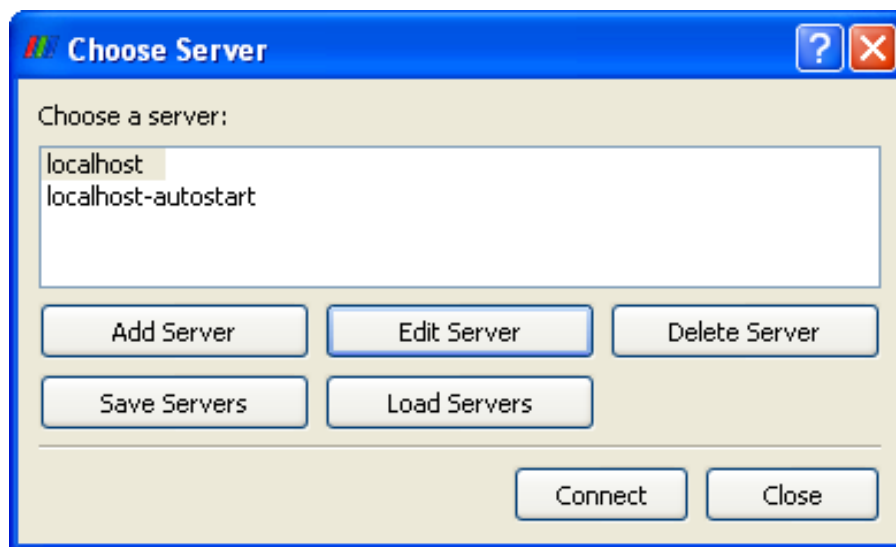


Figure 1. The Choose Server dialog establishes connections to configured servers.

To configure a new server connection, click the **Add Server** button to add it to the list. The dialog box shown below will appear. Enter a name in the first text entry box; this is the name that will appear in the **Choose Server** dialog (shown above).

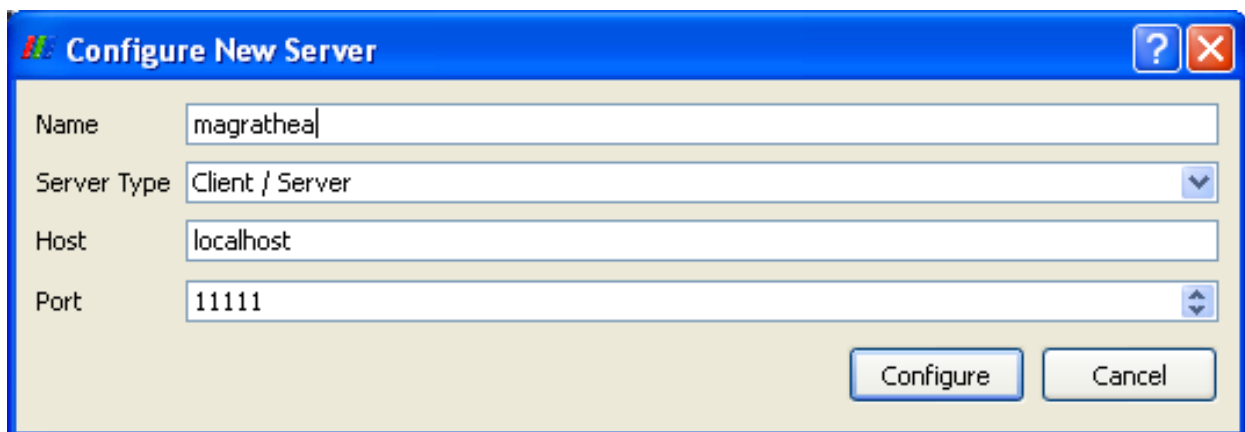


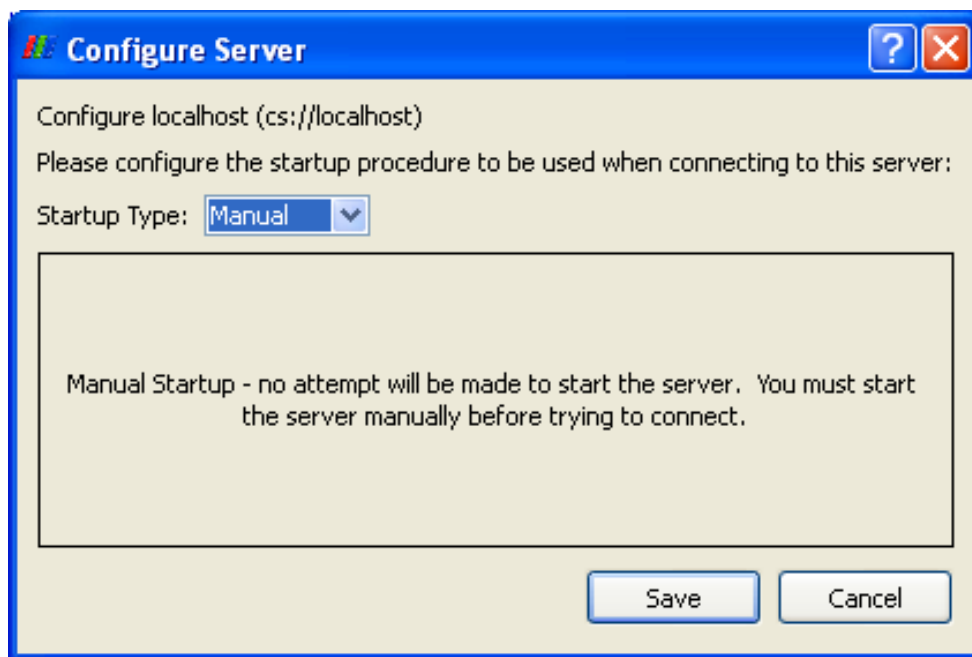
Figure 2. Configuring a new server.

Next select the type of connection you wish to establish from the **Server Type** menu. The possibilities are as follows. The “reverse connection” entries mean that the server connects to the client instead of the client connecting to the server. This may be necessary when the server is behind a firewall. Servers are usually run with multiple processes and on a machine other than where the client is running.

- **Client / Server:** Attach the ParaView client to a server.
- **Client / Server (reverse connection):** Connect a server to the ParaView client.
- **Client / Data Server / Render Server:** Attach the ParaView client to separate data and render servers.
- **Client / Data Server / Render Server (reverse connection):** Attach both a data and a render server to the ParaView client.

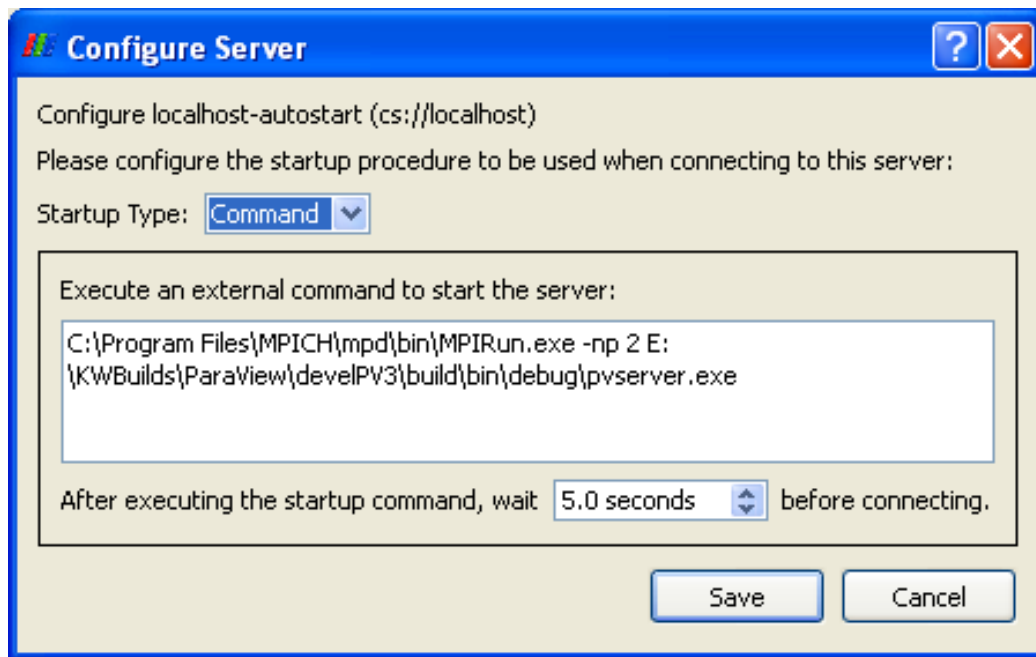
In either of the client / server modes, you must specify the name or IP address of the host machine (node 0) for the server. You may also enter a port number to use, or you may use the default (11111). If you are running in client / data server / render server mode, you must specify one host machine for the data server and another for the render server. You will also need two port numbers. The default one for the data server is 11111; the default for the render server is 22221.

When all of these values have been set, click the **Configure** button at the bottom of the dialog. This will cause the **Configure Server** dialog box, shown below, to appear. You must first specify the startup type. The options are **Command** and **Manual**. Choose **Manual** to connect to a server that has been started or will be started externally, on the command line for instance, outside of the ParaView user interface. After selecting **Manual**, click the **Save** button at the bottom of the dialog.



**Figure 3. Configure the server manually. It must be started outside of ParaView.**

If you choose the **Command** option, in the text window labeled **Execute an external command to start the server:**, you must give the command(s) and any arguments for starting the server. This includes commands to execute a command on a remote machine (e.g., `ssh`) and to run the server in parallel (e.g., `mpirun`). You may also specify an amount of time to wait after executing the startup command(s) and before making the connection between the client and the server(s). (See the spin box at the bottom of the dialog.) When you have finished, click the **Save** button at the bottom of the dialog.



**Figure 4. Enter a command that will launch a new server.**

Clicking the **Save** button in the **Configure Server** dialog will return you to the **Choose Server** dialog (shown earlier in this section). The server you just configured will now be in the list of servers you may choose. Thereafter, whenever you run ParaView you can connect to any of the servers that you have configured. You can also give the ParaView client the `--server=server_config_name` command-line argument to make it automatically connect to any of the servers from the list when it starts.

You can save and/or load server configurations to and/or from a file using the **Save Servers** and **Load Servers** buttons, respectively, on the **Choose Server** dialog. This is how some visualization centers provide system wide server configurations to allow the novice user to simply click his or her choice and connect to an already configured ParaView server. The format of the XML file for saving the server configurations is discussed online at [http://paraview.org/Wiki/Server\\_Configuration](http://paraview.org/Wiki/Server_Configuration).

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# Parallel Rendering and Large Displays

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## About Parallel Rendering

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One of ParaView's strengths is its ability to offload the often demanding rendering task. By offload, we mean that ParaView allows you to connect to a remote machine, ideally one that is closer to the data and to high-end rendering hardware, do the rendering on that machine and still interact with the data from a convenient location, such as the desktop or laptop in your office.

Abstracting away the location where rendering takes place opens up many possibilities. First, it opens up the possibility to parallelize the job of rendering to make it possible to render huge data sets at interactive rates. Rendering is done in the parallel Render Server component, which may be part of or separate from the parallel Data Server component. In the next section we describe how parallel rendering works and explain the controls you have over it. Second, huge datasets often require high resolution displays to view the intricate details within while maintaining a high level view to maintain context. In the following section we explain how ParaView can be used to drive tile display walls. Lastly, with the number of displays free to vary, it becomes possible to use ParaView to drive multi-display Virtual Reality systems. That is described in the final section of this chapter.

## Parallel Rendering

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### X Forwarding - Not a good idea

Parallel Rendering implies that many processors have some context to render their pixels into. Even though X11 forwarding might be available, **you should not run the client remotely and forward its X calls**. ParaView will be far more efficient if you let ParaView directly handle the data transfer between local and remote machines. When doing hardware accelerated rendering in GPUs this implies having X11 or Windows display contexts (either offscreen or onscreen). Otherwise this implies using offscreen Mesa (OSMesa) linked in to ParaView to do the rendering entirely offscreen into software buffers.

### Onscreen GPU Accelerated rendering via X11 Connections

One of the most common problems people have with setting up the ParaView server is allowing the server processes to open windows on the graphics card on each process's node. When ParaView needs to do parallel rendering, each process will create a window that it will use to render. This window is necessary because you need the X window before you can create an OpenGL context on the graphics hardware.

There is a way around this. If you are using the Mesa as your OpenGL implementation, then you can also use the supplemental OSMesa library to create an OpenGL context without an X window. However, Mesa is strictly a CPU rendering library so, **use the OSMesa solution if and only if your server hardware does not have rendering hardware**. If your cluster does not have graphics hardware, then compile ParaView with OSMesa support and use the `--use-offscreen-rendering` flag when launching the server.

Assuming that your cluster does have graphics hardware, you will need to establish the following three things.

1. Have xdm run on each cluster node at startup. Although xdm is almost always run at startup on workstation installations, it is not as commonplace to be run on cluster nodes. Talk to your system administrators for help in setting this up.
-

2. Disable all security on the X server. That is, allow any process to open a window on the x server without having to log in. Again, talk to your system administrators for help.
3. Use the `-display` flag for `pvserver` to make sure that each process is connecting to the display `localhost:0` (or just `:0`).

To enable the last condition, you would run something like

```
mpirun -np 4 ./pvserver -display localhost:0
```

An easy way to test your setup is to use the `glxgears` program. Unlike `pvserver`, it will quickly tell you (or, rather, fail to start) if it cannot connect to the local X server.

```
mpirun -np 4 /usr/X11R6/bin/glxgears -display localhost:0
```

## Offscreen software rendering via OSMesa

When running ParaView in a parallel mode, it may be helpful for the remote rendering processes to do their rendering in offscreen buffers. For example, other windows may be displayed on the node(s) where you are rendering; if these windows cover part of the rendering window, depending on the platform and graphics capabilities they might even be captured as part of the display results from that node. A similar situation could occur if more than one rendering process is assigned to a single machine, and the processes share a display. Also, in some cases the remote rendering nodes are not directly connected to a display and otherwise if your cluster does not have graphics hardware, then compile ParaView with OSMesa support and use the `--use-offscreen-rendering` flag when launching the server.

The first step to compiling OSMesa support is to make sure that you are compiling with the Mesa 3D Graphics Library <sup>[1]</sup>. It is difficult to tell an installation of Mesa from any other OpenGL implementation (although the existence of an `osmesa.h` header and a `libOSMesa` library is a good clue). If you are not sure, you can always download your own copy from <http://mesa3d.org>. We recommend using either Mesa version 7.6.1 or 7.9.1.

There are three different ways to use Mesa as ParaView's OpenGL library:

- You can use it purely as a substitute for GPU enabled onscreen rendering. To do this, set the CMake variable **OPENGL\_INCLUDE\_DIR** to point to the Mesa include directory (the one containing the GL subdirectory), and set the **OPENGL\_gl\_LIBRARY** and **OPENGL\_glu\_LIBRARY** to the libGL and libGLU library files, respectively.

| Variable              | Value              | Description   |
|-----------------------|--------------------|---|
| PARAVIEW_BUILD_QT_GUI | ON                 |   |
| VTK_USE_COCOA         | ON                 | Mac Only. X11 is not supported.                     |
| VTK_OPENGL_HAS_OSMESA | OFF                | Disable off screen rendering.                       |
| OPENGL_INCLUDE_DIR    | <mesa include dir> | Set this to the include directory in MESA.          |
| OPENGL_gl_LIBRARY     | libGL              | Set this to the libGL.a or libGL.so file in MESA.   |
| OPENGL_glu_LIBRARY    | libGLU             | Set this to the libGLU.a or libGLU.so file in MESA. |

- You can use it as a supplement to onscreen rendering. This mode requires that you have a display (X11 is running). In addition to specifying the GL library (which may be a GPU implementation of the Mesa one above), you must tell ParaView where Mesa's OSMesa library is. Do that by turning the **VTK\_OPENGL\_HAS\_OSMESA** variable to ON. After you configure again you will see a new CMake variable called **OSMESA\_LIBRARY**. Set this to the libOSMesa library file.

| Variable              | Value              | Description                                       |
|-----------------------|--------------------|---|
| PARAVIEW_BUILD_QT_GUI | ON                 |   |
| VTK_USE_COCOA         | ON                 | Mac Only. X11 is not supported.                   |
| VTK_OPENGL_HAS_OSMESA | ON                 | Turn this to ON to enable software rendering.     |
| OSMESA_INCLUDE_DIR    | <mesa include dir> | Set this to the include directory for MESA.       |
| OPENGL_INCLUDE_DIR    | <mesa include dir> | Set this to the include directory for MESA.       |
| OPENGL_gl_LIBRARY     | libGL              | Set this to the libGL.a or libGL.so file.         |
| OPENGL_glu_LIBRARY    | libGLU             | Set this to the libGLU.a or libGLU.so file.       |
| OSMESA_LIBRARY        | libOSMesa          | Set this to the libOSMesa.a or libOSMesa.so file. |

- You can use it for pure offscreen rendering, which is necessary when there is no display. To do this, make sure that the **OPENGL\_gl\_LIBRARY** variable is empty and that **VTK\_USE\_X** is off. Specify the location of OSMesa and **OPENGL\_glu\_LIBRARY** as above and turn on the **VTK\_USE\_OFFSCREEN** variable.

| Variable              | Value              | Description                                       |
|-----------------------|--------------------|---|
| PARAVIEW_BUILD_QT_GUI | OFF                | When using offscreen rendering there is no gui    |
| VTK_USE_COCOA         | OFF                | Mac only.   |
| VTK_OPENGL_HAS_OSMESA | ON                 | Turn this to ON to enable Off Screen MESA.        |
| OSMESA_INCLUDE_DIR    | <mesa include dir> | Set this to the include directory for MESA.       |
| OPENGL_INCLUDE_DIR    | <mesa include dir> | Set this to the include directory for MESA.       |
| OPENGL_gl_LIBRARY     | <empty>            | Set this to empty.                                |
| OPENGL_glu_LIBRARY    | libGLU             | Set this to the libGLU.a or libGLU.so file.       |
| OSMESA_LIBRARY        | libOSMesa          | Set this to the libOSMesa.a or libOSMesa.so file. |

Once again, once you build with OSMesa support, it will not take effect unless you launch the server with the `--use-offscreen-rendering` flag or alternatively, set the **PV\_OFFSCREEN** environment variable on the server to 1.

## Compositing

Given that you are connected to a server which is capable of rendering, you have a choice of whether to do the rendering remotely or to do it locally. ParaView's server performs all data-processing tasks. This includes generation of a polygonal representation of the full data set and of decimated LOD models. Once the data is generated on the server, it is sometimes better to do the rendering remotely and ship pixels to the client for display, and sometimes better to instead shift geometry to the client and have the client render it locally.

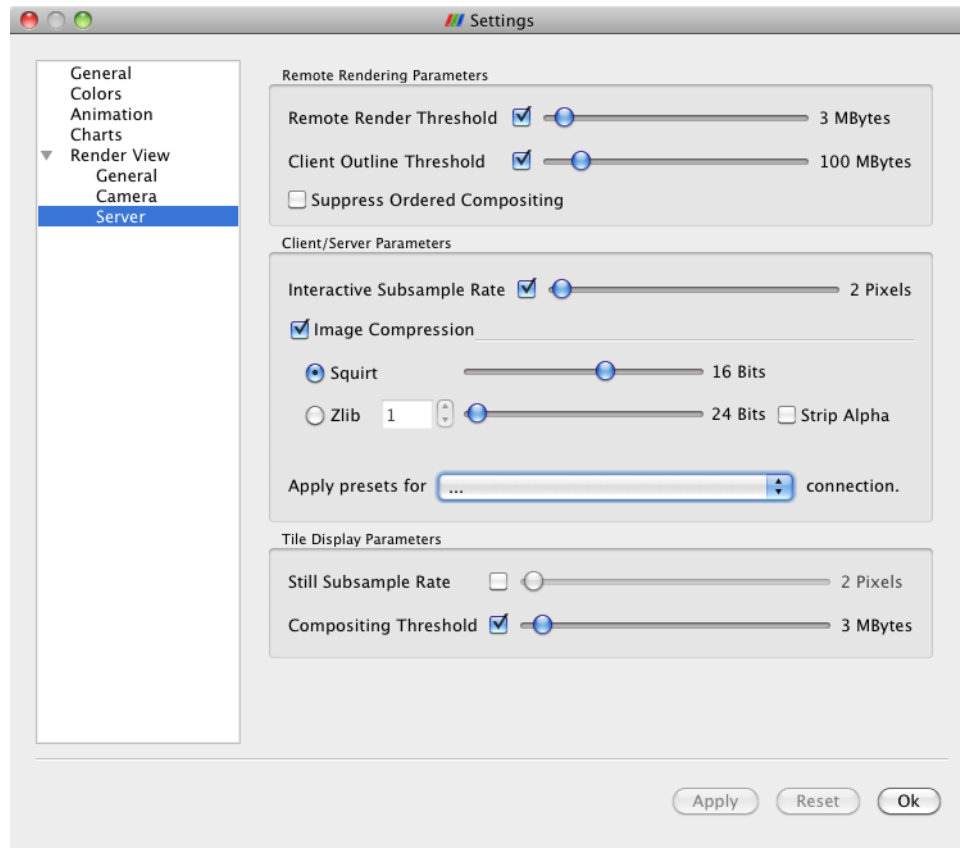
In many cases, the polygonal representation of the data set is much smaller than the original data set. (In an extreme case, a simple outline may be used to represent a very large structured mesh.) In these cases, it may be better to transmit the polygonal representation from the server to the client, and then let the client render it. The client can render the data repeatedly, when the viewpoint is changed for instance, without causing additional network traffic. Only when the data changes will network traffic occur. If the client workstation has high-performance rendering hardware, it can sometimes render even large data sets interactively in this way.

The second option is to have each node of the server render its geometry and send the resulting images to the client for display. There is a penalty per rendered frame for compositing images and sending the image across the network. However, ParaView's image compositing and delivery is very fast, and there are many options to ensure interactive rendering in this mode. Therefore, although small models may be collected and rendered on the client interactively,



ParaView's distributed rendering can render models of all sizes interactively.

ParaView automatically chooses a rendering strategy to achieve the best rendering performance. You can control the rendering strategy explicitly, forcing rendering to occur entirely on the server or entirely on the client for example, by choosing **Settings...** from the **Edit** menu of ParaView. Double click on **Render View** from the window on the left-hand side of the **Settings** dialog, and then click on **Server**. The rendering strategy parameters shown in Figure 10 will now be visible. Here we explain in detail the most important of these controls. For an explanation of all controls, see the Appendix.



**Figure 10. Parallel rendering parameters**

**Remote Render Threshold:** This slider determines how large the data set must be in order for parallel rendering with image compositing and delivery to be used (as opposed to collecting the geometry to the client). The value of this slider is measured in megabytes. Only when the entire data set consumes more memory than this value will compositing of images occur. If the check box beside the **Remote Render Threshold** slider is unmarked, then compositing will not happen; the geometry will always be collected. This is only a reasonable option when you can be sure the data set you are using is very small. In general, it is safer to move the slider to the right than to uncheck the box.

ParaView uses IceT to perform image compositing. IceT is a parallel rendering library that takes multiple images rendered from different portions of the geometry and combines them into a single image. IceT employs several image-compositing algorithms, all of which are designed to work well on distributed memory machines. Examples of two such image-compositing algorithms are depicted in Figure 11 and Figure 12. IceT will automatically choose a compositing algorithm based on the current workload and available computing resources.

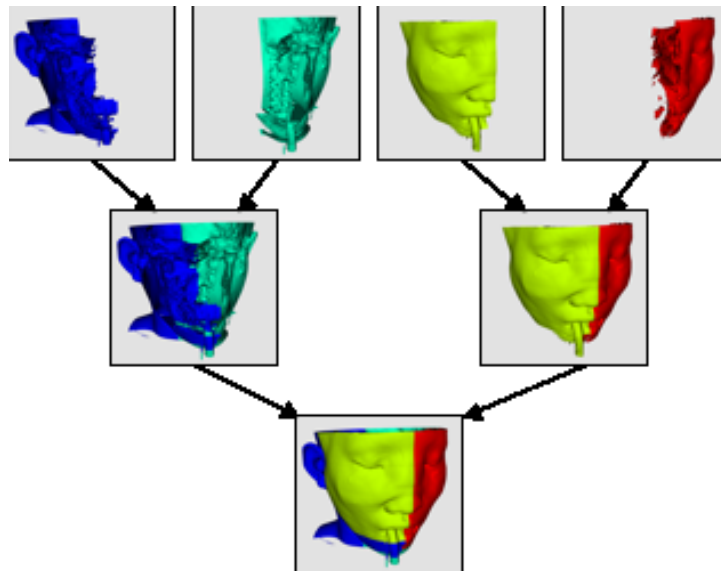


Figure 11. Tree compositing on four processes.

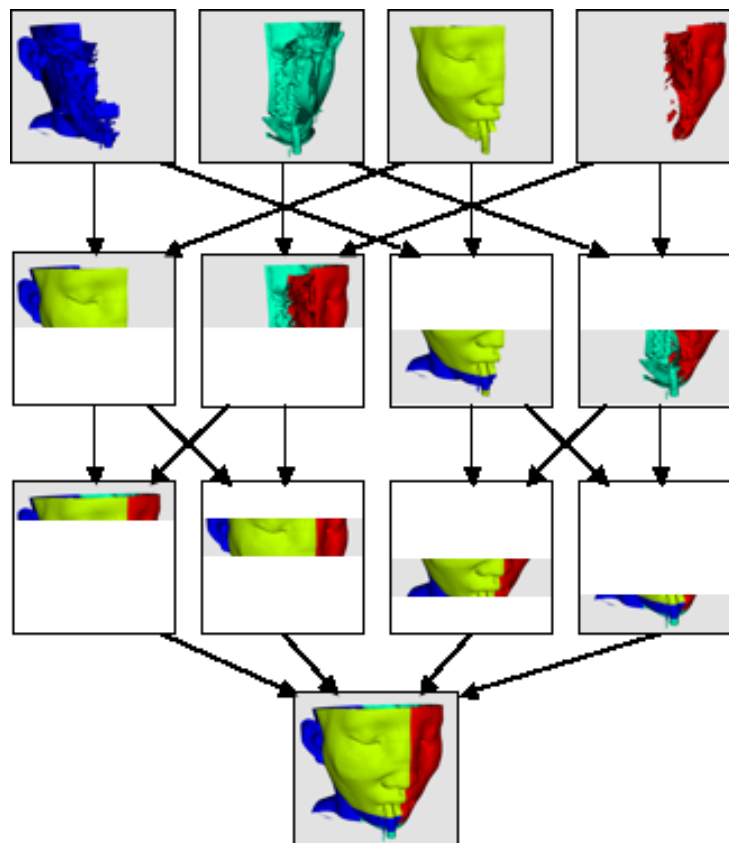


Figure 12. Binary swap on four processes.

**Interactive Subsample Rate:** The time it takes to composite and deliver images is directly proportional to the size of the images. The overhead of parallel rendering can be reduced by simply reducing the size of the images. ParaView has the ability to subsample images before they are composited and inflate them after they have been composited. The **Interactive Subsample Rate** slider specifies the amount by which images are subsampled. This is measured in pixels, and the subsampling is the same in both the horizontal and vertical directions. Thus a subsample rate of 2 will result in an image that is  $\frac{1}{4}$  the size of the original image. The image is scaled to full size before it is displayed on the user interface, so the higher the subsample rate, the more obviously pixilated the image will be during interaction as demonstrated in Figure 13. When the user is not interacting with the data, no subsampling will be used. If you want subsampling to always be off, unmark the check box beside the **Interactive Subsample Rate**

slider.

[[File:ParaView\_UsersGuide\_NoSubsampling.png link=]] [[File:ParaView\_UsersGuide\_TwoPixelSubsampling.png link=]] [[File:ParaView\_UsersGuide\_EightPixelSubsampling.png link=]]

No Subsampling                      Subsample Rate: 8 pixels  
    Subsample  
    Rate: 2  
    pixels

**Figure 13. The effect of subsampling on image quality**

**Squirt Compression:** When ParaView is run in client/server mode, ParaView uses image compression to optimize the image transfer. The compression uses an encoding algorithm optimized for images called SQUIRT (developed at Sandia National Laboratories).

SQUIRT uses simple run-length encoding for its compression. A run-length image encoder will find sequences of pixels that are all the same color and encode them as a single run length (the count of pixels repeated) and the color value. ParaView represents colors as 24-bit values, but SQUIRT will optionally apply a bit mask to the colors before comparing them. Although information is lost when this mask is applied, the sizes of the run lengths are increased, and the compression gets better. The bit masks used by SQUIRT are carefully chosen to match the color sensitivity of the human visual system. A 19-bit mask employed by SQUIRT greatly improves compression with little or no noticeable image artifacts. Reducing the number of bits further can improve compression even more, but it can lead to more noticeable color-banding artifacts.

The **Squirt Compression** slider determines the bit mask used during interactive rendering (i.e., rendering that occurs while the user is changing the camera position or otherwise interacting with the data). During still rendering (when the user is not interacting with the data), lossless compression is always used. The check box to the left of the **Squirt Compression** slider toggles whether the SQUIRT compression algorithm is used at all.

## References

[1] <http://mesa3d.org>

# Tile Display Walls

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## Tiled Display

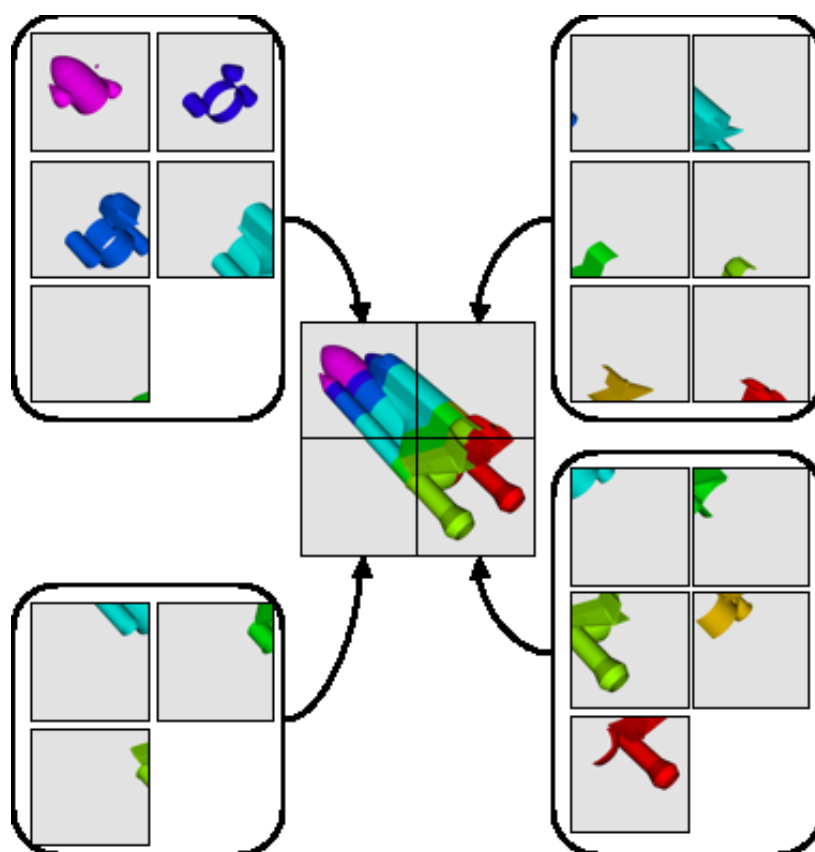
ParaView's parallel architecture makes it possible to visualize massive amounts of data interactively. When the data is of sufficient resolution that parallel processing is necessary for interactive display, it is often the case that high-resolution images are needed to inspect the data in adequate detail. If you have a 2D grid of display devices, you can run ParaView in tiled display mode to take advantage of it.

To put ParaView in tiled display mode, give `pvserver` (or `pvrenderserver`) the X and Y dimensions of the 2D grid with the `--tile-dimensions-x` (or `-tdx`) and `--tile-dimensions-y` (or `-tdy`) arguments. The X and Y dimensions default to 0, which disables tiled display mode. If you set only one of them to a positive value on the command line, the other will be set to 1. In tiled display mode, there must be at least as many server (in client / server mode) or render server (in client / data server / render server mode) nodes as tiles. The example below will create a 3 x 2 tiled display.

**`pvserver -tdx=3 -tdy=2`**

Unless you have a high end display wall, it is likely that each monitor's bezel creates a gap between the images shown in your tiled display. You can compensate for the bezel width by considering them to be like mullions on a window. To do that, specify the size of the gap (in pixels) through the `--tile-mullion-x` (or `-tmx`) and `--tile-mullion-y` (or `-tmy`) command line arguments.

The IceT library, which ParaView uses for its image compositing, has custom compositing algorithms that work on tiled displays. Although compositing images for large tiled displays is a compute-intensive process, IceT reduces the overall amount of work by employing custom compositing strategies and removing empty tiles from the computation as demonstrated in Figure 14. If the number of nodes is greater than the number of tiles, then the image compositing work will be divided amongst all the processes in the render server. In general, rendering to a tiled display will perform significantly better if there are many more nodes in the cluster than tiles in the display it drives. It also greatly helps if the geometry to be rendered is spatially distributed. Spatially distributed data is broken into contiguous pieces that are contained in small regions of space and are therefore rendered to smaller areas of the screen. IceT takes advantage of this property to reduce the amount of work required to composite the final images. ParaView includes the **D3** filter that redistributes data amongst the processors to ensure a spatially distributed geometry and thus improves tiled rendering performance.



**Figure 14. Compositing images for 8 processes on a 4-tile display.**

Unlike other parallel rendering modes, composited images are not delivered to the client. Instead, image compositing is reserved for generating images on the tiled display, and the desktop renders its own images from a lower resolution version of the geometry to display in the UI. In tiled display mode, ParaView automatically decimates the geometry and sends it to the client to make this happen. However, when the data is very large, even a decimated version of the geometry can overwhelm the client. In this case, ParaView will replace the geometry on the client with a bounding box.

You have several controls at run time over the tiled rendering algorithm that you can tune to maintain interactivity while visualizing very large data on very high resolution tiled displays. These are located on the **Tile Display Parameters** section of the **Render View / Server** page of the application settings dialog. These controls are described in the Application Settings section of the Appendix.

# CAVE Displays

## Introduction

- ParaView has basic support for visualization in CAVE like environments. These environments consists of multiple displays which are configured based on some coordinates in the room. Typically the base coordiante of the room is the reference coordinate of the head-tracker.
- ParaView can operate in a Client-Server fashion. The CAVE modules leverages this by assigning different process on the server to manage different displays. The client acts as a central point of control, sending tracking values to each of the processes on the server.
- ParaView has a VRPN (Virtual Reality Private Network) client plugin which reads tracked information from any VRPN server and relays the value to each process managing the display.
- Displays are configured using the PVX configuration XML file.

## Prerequisites

- VRPN needs to be installed on the client side.
- MPI needs to be installed on the server side.
- Make sure to enable BUILD\_SHARED\_LIB, PARAVIEW\_USE\_MPI and PARAVIEW\_BUILD\_PLUGIN\_VRPNPlugin when building for CAVE support.
- Consequently VRPN\_INCLUDE\_DIR and VRPN\_LIBRARY should also be set.

## Structure of config file

```
<?xml version="1.0" ?>
<pvx>
  <Process Type="server|dataserver|renderserver">
    <!--
      The only supported Type values are "server", "dataserver" or
      "renderserver".
      This controls which executable this configuration is applicable
      to.
      There can be multiple <Process /> elements in the same pvx file.
      -----
      | Executable      | Applicable Process Type          |
      | pvserver       | server, dataserver, renderserver |
      | pvrenderserver | server, renderserver             |
      | pvdataserver   | server, dataserver              |
      |-----|
    -->
    <Machine name="hostname"
      Environment="DISPLAY=m1:0"
      LowerLeft="-1.0 -1.0 -1.0"
      LowerRight="1.0 -1.0 -1.0"
      UpperRight="1.0 1.0 -1.0">
    <!--
      There can be multiple <Machine> elements in a <Process> element,
```

```

    each one identifying the configuration for a process.
    All attributes are optional.
    name="hostname"
    Environment: the environment for the process.
    LowerLeft|LowerRight|UpperRight
-->
</Machine>
</Process>
</pvx>

```

## Example Config

- The following example is for a six sided cave with origin at (0,0,0).

```

<?xml version="1.0" ?>
<pvx>
  <Process Type="client" />
  <Process Type="server">
    <Machine Name="Front"
      Environment="DISPLAY=:0"
      LowerLeft=" -1 -1 -1"
      LowerRight=" 1 -1 -1"
      UpperRight=" 1  1 -1" />
    <Machine Name="Right"
      Environment="DISPLAY=:0"
      LowerLeft="  1 -1 -1"
      LowerRight=" 1 -1  1"
      UpperRight=" 1  1  1" />
    <Machine Name="Left"
      Environment="DISPLAY=:0"
      LowerLeft=" -1 -1  1"
      LowerRight="-1 -1 -1"
      UpperRight="-1  1 -1"/>
    <Machine Name="Top"
      Environment="DISPLAY=:0"
      LowerLeft=" -1  1 -1"
      LowerRight=" 1  1 -1"
      UpperRight=" 1  1  1"/>
    <Machine Name="Bottom"
      Environment="DISPLAY=:0"
      LowerLeft=" -1 -1  1"
      LowerRight=" 1 -1  1"
      UpperRight=" 1 -1 -1"/>
    <Machine Name="Back"
      Environment="DISPLAY=:0"
      LowerLeft="  1 -1  1"
      LowerRight="-1 -1  1"
      UpperRight="-1  1  1"/>
  </Process>

```

</pvx>

- A sample PVX is provided in ParaView/Documentation/cave.pvx. This can be used to play with different display configurations.

### Things to keep in mind

- pvx file should be specified as the last command line argument for any of the server processes.
- Typically the pvx file is specified for all the executables whose environment is being changed using the pvx file. In case of data-server/render-server configuration, if you are setting up the environment for the two processes groups, then the pvx file must be passed as a command line option to both the executables: pvdataserver and pvrenderserver.
- When running in parallel the file is read on all nodes, hence must be present on all nodes.

### Basic steps for a successful setup

- On a terminal run the server (6 processes each driving one display in the cave. Change the number of process according to the number of displays).

```
# PV_ICET_WINDOW_BORDERS=1 mpiexec -np 6 ./pvserver /path/to/ConfigFile.pvx
```

Note: PV\_ICET\_WINDOW\_BORDERS=1 disables the full screen and instead opens up a 400x400 window. Remove this environment variable to work in full screen mode

- Open another terminal to run the client. (note: this client connects to the server which open up 6 windows according to the config given in cave.pvx). Change the stereo-type according to preference (see # ./paraview --help).

```
# ./paraview --stereo --stereo-type=Anaglyph --server=localhost --vrpn --vrpn-address=Tracker0@localhost
```

- Connect the client to the server.
  - Enable the VRPN Plugin on client.
-



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# Scripted Control

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## Interpreted ParaView

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The ParaView client provides an easy to use GUI in which to visualize data with the standard window, menu, button controls. Driving ParaView with a mouse is intuitive but not easily reproducible and exact as is required for repetitive analysis and scientific result reproducibility. For this type of work it is much better to use ParaView's scripted interface. This is an alternative control path, that works alongside of or as a replacement for the GUI.

ParaView's native scripted interface uses the Python programming language. The interface allows one to programatically control ParaView's back end data processing and rendering engines. Note that this level of control is distinct from the Python Programmable Filter discussed in the Quantitative Analysis Chapter. That is a white box filter, that runs in parallel within the server and has direct access to each individual element of the data. Here we are discussing client side python control, where you control the entire pipeline, which may or may not include Python Programmable Filters.

In this chapter we discuss first ParaView's python scripted interface, which follows the same idioms as and thus be operated similarly to the way one works, in the GUI. Next we discuss the ParaView's python tools, which make it even easier to use ParaView python scripts including the ability to record script files from GUI actions and to augment the GUI with them. Lastly we discuss batch processing with ParaView.

## Python Scripting

---

NOTE: This document is based on ParaView 3.6 or higher. If you are using 3.4, go to the history page and select the version from May 13, 2009.

### ParaView and Python

ParaView offers rich scripting support through Python. This support is available as part of the ParaView client (paraview), an MPI-enabled batch application (pvbatch), the ParaView python client (pvpython) or any other Python-enabled application. Using Python, users and developers can gain access to the ParaView engine called Server Manager.

*Note: Server Manager is a library that is designed to make it easy to build distributed client-server applications.*

This document is a short introduction to ParaView's Python interface. You may also visit the Python recipes page for some examples.

### Quick Start - a Tutorial

#### Getting Started

To start interacting with the Server Manager, you have to load the "simple" module. This module can be loaded from any python interpreter as long as the necessary files are in PYTHONPATH. These files are the shared libraries located in the paraview binary directory and python modules in the paraview directory: paraview/simple.py, paraview/vtk.py etc. You can also use either pvpython (for stand-alone or client/server execution), pvbatch (for

---

non-interactive, distributed batch processing) or the python shell invoked from Tools -> Python Shell using the ParaView client to execute Python scripts. You do not have to set PYTHONPATH when using these. In this tutorial, I will be using the python integrated development environment IDLE. My PYTHONPATH is set to the following.

```
/Users/berk/work/paraview3-build/bin:/Users/berk/work/paraview3-build/Utilities/VTKPythonWrapping/site-packages
```

Note: This for older versions of ParaView this was  
 /Users/berk/work/paraview3-build/bin:/Users/berk/work/paraview3-build/Utilities/VTKPythonWrapping  
 --Andy.bauer 23 July 2010.

You may also need to set your path variable for searching for shared libraries (i.e. PATH on Windows and LD\_LIBRARY\_PATH on Unix/Linux/Mac). The corresponding LD\_LIBRARY\_PATH would be:

```
/Users/berk/work/paraview3-build/bin
```

(Under WindowsXP for a debug build of paraview, I set BOTH my PATH and PYTHONPATH environment variables to include \${BUILD}/bin/Debug and \${BUILD}/Utilities/VTKPythonWrapping to make it work.  
 --DaveDemarle 21:09, 29 June 2009 (UTC))

This is on my Mac and using the build tree. In IDLE, let's start by loading the servermanager module.

```
>>> from paraview.simple import *
```

Note: Importing the paraview module directly is deprecated although still possible for backwards compatibility. This document refers to the simple module alone.

In this example, we will use ParaView in the stand-alone mode. Connecting to a ParaView server running on a cluster is covered later in this document.

## Creating a Pipeline

The simple module contains many functions to instantiate sources, filters and other related objects. You can get a list of object this module can create from ParaView's online help (from help menu or here: <http://paraview.org/OnlineHelpCurrent/>)

Let's start by creating a Cone object:

```
>>> cone = Cone()
```

You can get some documentation about the cone object using help().

```
>>> help(cone)
Help on Cone in module paraview.servermanager object:

class Cone(SourceProxy)
| The Cone source can be used to add a polygonal cone to the 3D
| scene. The output of the
| Cone source is polygonal data.
|
| Method resolution order:
|   Cone
|   SourceProxy
|   Proxy
|   __builtin__.object
|
| Methods defined here:
```

```
|
| Initialize = aInitialize(self, connection=None)
|
|
|-----
| Data descriptors defined here:
|
| Capping
|     If this property is set to 1, the base of the cone will be
capped with a filled polygon.
Otherwise, the base of the cone will be open.
|
| Center
|     This property specifies the center of the cone.
|
| Direction
|     Set the orientation vector of the cone. The vector does not
have to be normalized. The cone
will point in the direction specified.
|
| Height
|     This property specifies the height of the cone.
|
| Radius
|     This property specifies the radius of the base of the cone.
|
| Resolution
|     This property indicates the number of divisions around the
cone. The higher this number, the
closer the polygonal approximation will come to representing a cone,
and the more polygons it will
contain.
|
|...
```

This gives you a full list of properties. Let's check what the resolution property is set to.

```
>>> cone.Resolution
6
```

You can increase the resolution as shown below.

```
>>> cone.Resolution = 32
```

Alternatively, we could have specified a value for resolution when creating the object.

```
>>> cone = Cone(Resolution=32)
```

You can assign values to any number of properties during construction using keyword arguments. Let's also change the center.

```
>>> cone.Center
[0.0, 0.0, 0.0]
>>> cone.Center = [1, 2, 3]
```

Vector properties such as this one support setting and getting of individual elements as well as slices (ranges of elements).

```
>>> cone.Center[0:2] = [2, 4]
>>> cone.Center
[2.0, 4.0, 3.0]
```

Next, let's apply a shrink filter to the cone:

```
>>> shrinkFilter = Shrink(cone)
>>> shrinkFilter.Input
<paraview.servermanager.Cone object at 0xaf701f0>
```

At this point, if you are interested in getting some information about the output of the shrink filter, you can force it to update (which will also cause the execution of the cone source). For details about VTK's demand-driven pipeline model used by ParaView, see one of the VTK books.

```
>>> shrinkFilter.UpdatePipeline()
>>> shrinkFilter.GetDataInformation().GetNumberOfCells()
33L
>>> shrinkFilter.GetDataInformation().GetNumberOfPoints()
128L
```

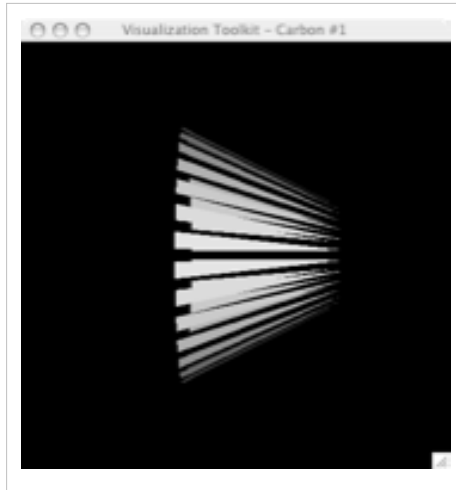
We will cover the `DataInformation` class in more detail later.

## Rendering

Now that we created a small pipeline, let's render the result. You will need two objects to render the output of an algorithm in a scene: a representation and a view. A representation is responsible for taking a data object and rendering it in a view. A view is responsible for managing a render context and a collection of representations. Simple creates a view by default. The representation object is created automatically with `Show()`.

```
>>> Show(shrinkFilter)
>>> Render()
```

Et voila:



In the example above, we assigned the value returned by `Cone()` and `Shrink()` to Python variables and used them to build our pipeline. ParaView keeps track of the last pipeline object created by the user. This allows us to accomplish everything we did above using the following code.

```
>>> from paraview.simple import *
# Create a cone and assign it as the active object
>>> Cone()
<paraview.servermanager.Cone object at 0x2910f0>
# Set a property of the active object
>>> SetProperties(Resolution=32)
# Apply the shrink filter to the active object
# Shrink is now active
>>> Shrink()
<paraview.servermanager.Shrink object at 0xaf64050>
# Show shrink
>>> Show()
<paraview.servermanager.UnstructuredGridRepresentation object at 0xaf57f90>
# Render the active view
>>> Render()
<paraview.servermanager.RenderView object at 0xaf57ff0>
```

This was a quick introduction to the `paraview.simple` module. In the following sections, we will discuss the Python interface in more detail and introduce more advanced concepts.

## paraview.simple Module

The simple module is a ParaView component written using Python on top of the Server Manager C++ library. Its purpose is to make it easier to create ParaView data analysis and visualization pipelines using Python. The simple module can be loaded from Python interpreters running in several applications.

- `pvpython`: The `pvpython` application, distributed with the ParaView application suite, is a Python client to the ParaView servers. It supports interactive execution as well as batch execution.
- `pvbatch`: The `pvbatch` application, also distributed with the ParaView application suite, is a Python application designed to run batch scripts on distributed servers. When ParaView is compiled with MPI, `pvbatch` can be launched as an MPI program. In this mode, the first node will load a Python script specified as a command-line argument and execute it using a special built-in connection on all nodes. This application does not support

interactive execution.

- paraview: Python scripts can be run from the paraview client using the Python shell that is invoked from Tools -> Python Shell. The Python shell supports interactive mode as well as loading of scripts from file.
- External Python interpreter: Any Python-capable application can load the paraview.simple module if the right environment is configured. For this to work, you either have to install the paraview Python modules (including the right shared libraries) somewhere in sys.path or you have to set PYTHONPATH to point to the right locations.

## Overview

The paraview.simple module contains several Python classes designed to be Python-friendly as well as all classes wrapped from the C++ Server Manager library. The following sections cover the usage of this module and occasionally the paraview.servermanager module which is lower level.

## Connecting to a Server

ParaView can run in two modes: stand-alone and client/server where the server is usually a visualization cluster. In this section, we discuss how to establish a connection to a server when using ParaView in the client/server mode. If you are using the ParaView graphical interface, you should use Connect from the File menu to connect to a server. If you are using ParaView from a Python shell (not the Python console that is part of the graphical interface), you need to use servermanager.Connect() to connect a server. *Note: you cannot connect to the ParaView application from a stand-alone Python shell. You can only connect to a server.* This method takes 4 arguments, all of which have default values.

```
def Connect(ds_host=None, ds_port=11111, rs_host=None, rs_port=11111)
```

When connecting to a server (pvserver), specify only the first 2 arguments. These are the server name (or IP address) and port number.

When connecting to a data-server/render-server pair, you have to specify all four arguments. The first 2 are the host name (or IP address) and port number of the data server, the last 2 those of the render server. Here are some examples.

```
# Connect to pvserver running on amber1 (first node of our test cluster)
# using the default port 11111
>>> Connect('amber1')

# Connect to pvdataserver running on the amber cluster, pvrenderserver
# running on Berk's desktop
>>> Connect('amber1', 12000, 'kamino', 11111)
```

Note: Connect() will return None on failure. To be safe, you should check the return value of Connect().

## Getting Help

You can access the documentation of all Proxy types by using Python's built-in help.

```
>>> help(paraview.simple.Cone)
Help on function CreateObject in module paraview.simple:

CreateObject(*input, **params)

    The Cone source can be used to add a polygonal cone to the 3D
    scene. The output of the
```

```
Cone source is polygonal data.
```

To get the full documentation, you have to create an instance.

```
>>> c = Cone()
>>> help(c)
```

This documentation is automatically generated from the Server Manager configuration files and is identical to the class documentation found under the paraview Help menu as well as here <http://paraview.org/OnlineHelpCurrent/>. Beyond this document and the online help, there are a few useful documentation sources.

- The ParaView Guide: <http://www.kitware.com/products/paraviewguide.html>
- The ParaView Wiki: <http://paraview.org/Wiki/ParaView>
- The ParaView source documentation: <http://www.paraview.org/doc/>

If you are interested in learning more about the Visualization Toolkit that is at the foundation of ParaView, visit <http://vtk.org>.

## Proxies and Properties

### Proxies

The VTK Server Manager design uses the Proxy design pattern (*See Design Patterns: Elements of Reusable Object-Oriented Software by Erich Gamma, Richard Helm, Ralph Johnson and John Vlissides for details*). Quoting from Wikipedia: “A proxy, in its most general form, is a class functioning as an interface to another thing. The other thing could be anything: a network connection, a large object in memory, a file, or some other resource that is expensive or impossible to duplicate”. In the case of Server Manager, a Proxy object acts as a proxy to one or more VTK objects. Most of the time, these are server-side objects and are distributed to the server nodes. Proxy objects allow you to interact with these object as if you directly have access to them, manipulate them and obtain information about them. When creating visualization pipelines, you create proxies instead of VTK objects.

```
>>> sphereSource = vtk.vtkSphereSource() # VTK-Python script

>>> sphereSourceP = Sphere() # ParaView script
```

A proxy also provides an interface to modify the properties of the objects it maintains. For example, instead of

```
>>> sphereSource.SetCenter(1.0, 1.0, 0.0)
```

you can write the following.

```
>>> sphere.Center = [1.0, 1.0, 0.0]
```

When a pipeline object proxy is created, it is set as the active object. You can also set an object as the active one. This is equivalent to clicking on an object in the pipeline browser.

```
>>> c = Cone()
<paraview.servermanager.Cone object at 0xaf73090>
>>> GetActiveSource()
<paraview.servermanager.Cone object at 0xaf73090>
>>> Shrink()
<paraview.servermanager.Shrink object at 0xb4f8610>
# Make the cone active
```

```
>>> SetActiveSource(c)
```

When dealing with objects created through the graphical interface or by loading a state, it is useful to be able to search through existing pipeline objects. To accomplish this, you can use `GetSources()` and `FindSource()`. `GetSources()` returns a dictionary of (name, id), object pairs. Since multiple objects can have the same name, the (name,id) pair identifies objects uniquely. `FindSource()` returns an object given its name. If there are more than one objects with the same name, the first one is returned.

```
>>> Cone()
<paraview.servermanager.Cone object at 0xaf73090>
>>> GetActiveSource()
<paraview.servermanager.Cone object at 0xaf73090>
>>> Shrink()
<paraview.servermanager.Shrink object at 0xb4f8610>
>>> SetActiveSource(c)
```

To delete pipeline objects, you need to use the `Delete()` function. Simply letting a Python variable go out of scope is not enough to delete the object. Following the example above

```
# Delete the cone source
>>> Delete(c)
# To fully remove the cone from memory, get rid of the
# variable too
>>> del c
```

## Properties

Property objects are used to read and modify the properties of pipeline objects. Each proxy has a list of properties defined in the Server Manager configuration files. The property interface of the Server Manager C++ library is somewhat cumbersome. Here is how you can set the radius property of a sphere source.

```
>>> rp = sphere.GetProperty("Radius")
>>> rp.SetElement(0, 2)
1
>>> sphere.UpdateProperty("Radius")
```

The `servermanager` module makes property access much easier by defining Python property accessors for property objects:

```
>>> sphere.Radius = 3
```

Here `Radius` is a Python property which, when a value is assigned to it, calls `sphere.SetPropertyWithName("Radius",3)`. Properties can also be passed to the function creating the object:

```
>>> cone = Cone(Radius=0.5, Center=[1, 0.5, 0])
<paraview.servermanager.Cone object at 0xaf73090>
```

□ You can also use the `SetProperties()` function to set property values.

```
>>> SetProperties(cone, Radius=0.2, Center=[2, 0.5, 0])
```

If the first argument is not specified, the active object is used. You also use `SetDisplayProperties()` and `SetViewProperties()` to set display (representation) and view properties respectively.

All Property classes define the following methods.



- `__len__()`
- `__getitem__()`
- `__setitem__()`
- `__getslice__()`
- `__setslice__()`
- `GetData()`
- `SetData()`.

Therefore, all of the following are supported.

```
>>> sphere.Center
[0.0, 0.0, 0.0]
>>> sphere.Center[0] = 1
>>> sphere.Center[0:3] = [1,2,3]
>>> sphere.Center[0:3]
[1.0, 2.0, 3.0]
>>> len(sphere.Center)
3
```

ProxyProperty and InputProperty also define

- `append()`
- `__delitem__()`
- `__delslice__()`

to support `del()` and `append()`, similar to Python list objects.

VectorProperty is used for scalars, vectors and lists of integer and floating point numbers as well as strings. Most properties of this type are simple. Examples include `Sphere.Radius` (double scalar), `Sphere.Center` (vector of doubles), `a2DGlyph.Filled` (boolean), `a2DGlyph.GlyphType` (enumeration), `a3DText.Text` (string) and `Contour.Isosurfaces` (list of doubles). Some properties may be more complicated because they map to C++ methods with mixed argument types. Two good examples of this case are `Glyph.Scalars` and `ExodusIIReader.PointVariables`.

```
>>> reader = ExodusIIReader(FileName='.../can.ex2')
# These variables are currently selected
>>> reader.PointVariables
['DISPL', 'VEL', 'ACCL']
# These are available in the file
>>> reader.PointVariables.Available
['DISPL', 'VEL', 'ACCL']
# Enable the DISPL array only
>>> reader.PointVariables = ['DISPL']
# Force read
>>> reader.UpdatePipeline()
# Now check the output. Note: GlobalNodeId is generated automatically
by the reader.
>>> reader.PointData[:]
[Array: GlobalNodeId, Array: PedigreeNodeId, Array: DISPL]
```

This example demonstrates the use of `ExodusIIReader.PointVariables`. This is a `VectorProperty` that represents a list of array names. The underlying C++ function has a signature of `SetPointResultArrayStatus( const char* name, int flag )`. This method is usually called once per array to enable or disable it (i.e. to set whether the reader will read a

particular array).

Glyph.Scalars is a bit more complicated. This property allows the developer to select the scalar array with which to scale the glyphs.

```
>>> sph = Sphere()
>>> elev=Elevation(sph)
# Glyph the points of the sphere with spheres
>>> glyph=Glyph(elev, GlyphType='Sphere')
# Scale the glyph with the Elevation array
>>> glyph.Scalars = 'Elevation'
>>> glyph.Scalars
['POINTS', 'Elevation']
# The above shows the association of the array as well as its name.
# In this case, the array is associated with POINTS as it has to be
# since Glyph cannot scale by cell arrays. We could have done:
>>> glyph.Scalars = ['POINTS', 'Elevation']
# Enable scaling by scalars
>>> glyph.ScaleMode = 'scalar'
```

Here the property Scalars maps to SetInputArrayToProcess( int idx, int port, int connection, int fieldAssociation, const char \*name ) which has four integer arguments (some of which are enumeration) and 1 string argument (see *vtkAlgorithm documentation for details*).

Properties are either regular (push) or information (pull) properties. Information properties do not have a VTK method associated with them and are responsible for getting information from the server. A good example of an information property is TimestepValues which returns all time steps available in a file (if the reader supports time).

```
>>> reader = ExodusIIReader(FileName='../can.ex2')
>>> reader.TimestepValues
[0.0, 0.00010007373930420727, 0.00019990510190837085,
0.00029996439116075635, 0.00040008654468692839,
0.00049991923151537776, 0.00059993512695655227, 0.00070004, ...]
```

You can obtain a list of properties a proxy supports by using help(). However, this does not allow introspection programmatically. If you need to obtain information about a proxy's properties programmatically, you can use a property iterator:

```
>>> for prop in glyph:
    print type(prop), prop.GetXMLLabel()

<class 'paraview.servermanager.InputProperty'> Input
<class 'paraview.servermanager.VectorProperty'> Maximum Number of Points
<class 'paraview.servermanager.VectorProperty'> Random Mode
<class 'paraview.servermanager.ArraySelectionProperty'> Scalars
<class 'paraview.servermanager.ArraySelectionProperty'> Vectors
<class 'paraview.servermanager.VectorProperty'> Orient
<class 'paraview.servermanager.VectorProperty'> Set Scale Factor
<class 'paraview.servermanager EnumerationProperty'> Scale Mode
<class 'paraview.servermanager.InputProperty'> Glyph Type
<class 'paraview.servermanager.VectorProperty'> Mask Points
```

The XMLLabel is the text display by the graphical user interface. Note that there is a direct mapping from the XMLLabel to the property name. If you remove all spaces from the label, you get the property name. You can use the PropertyIterator object directly.

```
>>> it = iter(s)
>>> for i in it:
    print it.GetKey(), it.GetProperty()
```

## Domains

The Server Manager provides information about values that are valid for properties. The main use of this information is for the user interface to provide good ranges and choices in enumeration. However, some of this information is also very useful for introspection. For example, enumeration properties look like simple integer properties unless a (value, name) pair is associated with them. The Server Manager uses Domain objects to store this information. The contents of domains may be loaded from xml configuration files or computed automatically. Let's look at an example.

```
>>> s = Sphere()
>>> Show(s)
>>> dp = GetDisplayProperties(s)
>>> dp.Representation
'Surface'
# The current representation type is Surface. What other types
# are available?
>>> dp.GetProperty("Representation").Available
['Outline', 'Points', 'Wireframe', 'Surface', 'Surface With Edges']
# Choose outline
>>> dp.Representation = 'Outline'
```

## Source Proxies

Source proxies are proxies that represent pipeline objects (*For more information about VTK pipelines, see the VTK books: <http://vtk.org/buy-books.php>*). They have special properties to connect them as well as special method to query the meta-data of their output. To connect a source proxy to another, use one of its input properties.

```
# Either
>>> glyph = Glyph(elev)
# or
>>> glyph.Input = elev
```

The SourceProxy class provides several additional properties and methods that are specific to pipelines (See vtkSMSSourceProxy documentation for a full list).

- UpdatePipelineInformation(): This method calls UpdateInformation() on the VTK algorithm. It also calls UpdatePropertyInformation() to update any information properties.
- UpdatePipeline(): This method calls Update() on the VTK algorithm causing a pipeline execution if the pipeline changed. Another way of causing pipeline updates is to render. The render view updates all connected pipelines.
- GetDataInformation(): This method is used to obtain meta-data about one output. It is discussed further below.
- PointData and CellData properties discussed below.

There are two common ways of getting meta-data information from a proxy: information properties and DataInformation. Information properties are updated automatically every time UpdatePropertyInformation() and

UpdatePipelineInformation() are called. All you have to do is read the data from the property as usual. To get a DataInformation object from a source proxy use GetDataInformation(port=0). By default, this method returns data information for the first output. You can pass an optional port number to get information for another output. You can get detailed documentation on DataInformation by using help() and by reading online documentation for vtkPVDataInformation (<http://www.paraview.org/doc/nightly/html/classvtkPVDataInformation.html>). *Here are the use of some common methods.*

```
>>> di = glyph.GetDataInformation(0)
>>> di
<paraview.servermanager.DataInformation object at 0x2d0920d0>
>>> glyph.UpdatePipeline()
# Get the data type.
>>> di.GetDataClassName()
'vtkPolyData'
# Get information about point data.
>>> pdi = di.PointData
# We are now directly accessing the wrapper for a VTK class
>>> len(pdi)
1
# Get information for a point array
>>> ai = pdi[0]
>>> ai.GetRange(0)
(0.0, 0.5)
```

When meta-data is not enough and you need access to the raw data, you can use Fetch() to bring it to the client side. Note that this function is provided by the servermanager module. Fetch() has three modes:

- Append all of the data together and bring it to the client (only available for polygonal and unstructured datasets). Note: Do not do this if data is large otherwise the client will run out of memory.
- Bring data from a given process to the client.
- Use a reduction algorithm and bring its output to the client. For example, find the minimum value of an attribute.

Here is a demonstration.

```
>>> from paraview.simple import *
>>> Connect("kamino")
Connection (kamino:11111)
>>> s = Sphere()
# Get the whole sphere. DO NOT DO THIS IF THE DATA IS LARGE otherwise
# the client will run out of memory.
>>> allsphere = servermanager.Fetch(s)
getting appended
use append poly data filter
>>> allsphere.GetNumberOfPolys()
96
# Get the piece of the sphere on process 0.
>>> onesphere = servermanager.Fetch(s, 0)
getting node 0
>>> onesphere.GetNumberOfPolys()
48
```

```

# Apply the elevation filter so that we have a useful scalar array.
>>> elev = Elevation(s)
# We will use the MinMax algorithm to compute the minimum value of
# elevation. MinMax will be first applied on each processor. The results
# will then be gathered to the first node. MinMax will be then applied
# to the gathered results.
# We first create MinMax without an input.
>>> mm = MinMax(None)
# Set it to compute min
>>> mm.Operation = "MIN"
# Get the minimum
>>> mindata = servermanager.Fetch(elev, mm, mm)
applying operation
# The result is a vtkPolyData with one point
>>> mindata.GetPointData().GetNumberOfArrays()
2
>>> a0 = mindata.GetPointData().GetArray(1)
>>> a0.GetName()
'Elevation'
>>> a0.GetTuple1(0)
0.0

```

## Representations and Views

Once a pipeline is created, it can be rendered using representations and views. A view is essentially a “window” in which multiple representations can be displayed. When the view is a VTK view (such as `RenderView`), this corresponds to a collection of objects including `vtkRenderers` and a `vtkRenderWindow`. However, there is no requirement for a view to be a VTK view or to render anything. A representation is a collection of objects, usually a pipeline, that takes a data object, converts it to something that can be rendered and renders it. When the view is a VTK view, this corresponds to a collection of objects including geometry filters, level-of-detail algorithms, `vtkMappers` and `vtkActors`. The simple module automatically creates a view after connecting to a server (including the built-in connection when using the stand-alone mode). Furthermore, the simple module creates a representation the first time a pipeline object is displayed with `Show()`. It is easy to create new views.

```
>>> view = CreateRenderView()
```

`CreateRenderView()` is a special method that creates the render view appropriate for the `ActiveConnection` (or for another connection specified as an argument). It returns a sub-class of `Proxy`. Like the constructor of `Proxy`, it can take an arbitrary number of keyword arguments to set initial values for properties. Note that ParaView makes the view that was created last the active view. When using `Show()` without a view argument, the pipeline is shown in the active view. You can get a list of views as well as the active view as follows

```

>>> GetRenderViews()
[<paraview.servermanager.RenderView object at 0xaf64ef0>, <paraview.servermanager.RenderView object at 0xaf64b70>]
>>> GetActiveView()
<paraview.servermanager.RenderView object at 0xaf64b70>

```

You can also change the active view using `SetActiveView()`.

Once you have a render view, you can use pass it to `show` in order to select in which view a pipeline object is displayed. You can also pass it to `Render()` to select which view is rendered.

```
>>> Show(elev, GetRenderViews()[1])
<paraview.servermanager.GeometryRepresentation object at 0xaf64e30>
>>> Render(GetRenderViews()[1])
```

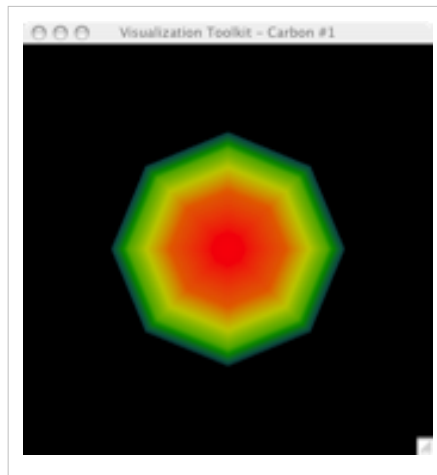
Notice that `Show()` returns a representation object (aka `DisplayProperties` in the `simple` module). This object can be used to manipulate how the pipeline object is displayed in the view. You can also access the display properties of an object using `GetDisplayProperties()`.

```
>>> dp = GetDisplayProperties(elev)
>>> dp
<paraview.servermanager.GeometryRepresentation object at 0xaf649d0>
```

Display properties and views have a large number of documented properties some of which are poorly documented. We will cover some them here.

```
>>> from paraview.simple import *
# Create a simple pipeline
>>> sph = Sphere()
>>> elev = Elevation(sph)
>>> Show(elev)
>>> Render()
# Set the representation type of elev
>>> dp = GetDisplayProperties(elev)
>>> dp.Representation = 'Points'
# Here is how you get the list of representation types
>>> dp.GetProperty("Representation").Available
['Outline', 'Points', 'Wireframe', 'Surface', 'Surface With Edges']
>>> Render()
# Change the representation to wireframe
>>> dp.Representation = 'Wireframe'
>>> Render()
# Let's get some information about the output of the elevation
# filter. We want to color the representation by one of it's
# arrays.
# Second array = Elevation. Interesting. Let's use this one.
>>> ai = elev.PointData[1]
>>> ai.GetName()
'Elevation'
# What is its range?
>>> ai.GetRange()
(0.0, 0.5)
# To color the representation by an array, we need to first create
# a lookup table. We use the range of the Elevation array
>>> dp.LookupTable = MakeBlueToRedLT(0, 0.5)
>>> dp.ColorAttributeType = 'POINT_DATA'
>>> dp.ColorArrayName = 'Elevation' # color by Elevation
>>> Render()
```

Here is the result: 



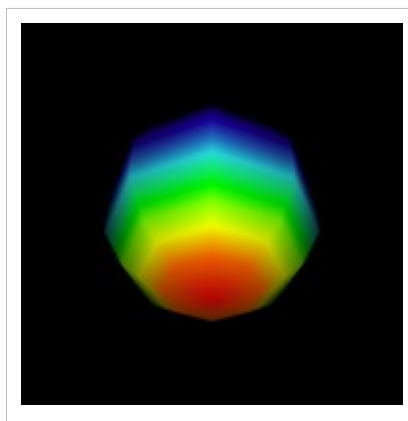
Once you create a scene, you will probably want to interact with the camera and `ResetCamera()` is likely to be insufficient. In this case, you can directly get the camera from the view and manipulate it. `GetActiveCamera()` returns a VTK object (not a proxy) with which you can interact.

```
>>> camera = GetActiveCamera()
>>> camera
<libvtkCommonPython.vtkCamera vtkobject at 0xe290>
>>> camera.Elevation(45)
>>> Render()
```

Another common thing to do is to save the view as an image. For this purpose, you can use the `WriteImage()` method provided by the view:

```
>> WriteImage("/Users/berk/image.png")
```

The resulting `image.png` looks like this. See the documentation for `WriteImage()` for details on choosing file type as well as a magnification factor to save images larger than the view size.



## Advanced Concepts

### Dealing with lookup tables

As shown earlier, you can use `MakeBlueToRedLt(min, max)` to create a lookup table. However this simply creates a new lookup table which the GUI won't be aware of. In ParaView Qt application, we have special lookup table management that ensures that the same lookup table is used for all arrays with same name and number of components. To reproduce the same behavior in Python, use `GetLookupTableForArray()`.

```
def GetLookupTableForArray(arrayname, num_components, **params):
    """Used to get an existing lookuptable for a array or to create one
    if none
    exists. Keyword arguments can be passed in to initialize the LUT if
    a new
    one is created. Returns the lookup table."""
    ....
```

This will create a new lookup table and associate it with that array, if none already exists. Any default arguments to be passed to the lookup table if a new one is created, can be specified as additional parameters. You can always change the properties on the lookup table returned by this function.

```
# To color the representation by an array, we need to first create
# a lookup table. We use the range of the Elevation array
>>> dp.LookupTable = GetLookupTableForArray("Elevation", 1,
                                             RGBPoints = [min, 0, 0, 1, max, 1, 0, 0],
                                             ColorSpace = "HSV")
>>> dp.ColorAttributeType = 'POINT_DATA'
>>> dp.ColorArrayName = 'Elevation' # color by Elevation
>>> Render()
```

### Loading State and Manipulating It

Let's say you created a complicated visualization using the paraview application and now you want to make slight changes to it and run it in a loop as a batch script. What do you do? The best way of dealing with this is to save your visualization state and then load it from Python. Let's say you have a state file saved as `myteststate.pvsm`:

```
>>> from paraview.simple import *
# Load the state
>>> servermanager.LoadState("/Users/berk/myteststate.pvsm")
# Make sure that the view in the state is the active one
>>> SetActiveView(GetRenderView())
# Now render
>>> Render()
# Get the list of sources
>>> GetSources()
{('Sphere1', '5'): <paraview.servermanager.Sphere object at 0xaf80e30>,
 ('Shrink1', '11'): <paraview.servermanager.Shrink object at 0xaf80df0>,
 ('Cone1', '8'): <paraview.servermanager.Cone object at 0xaf80cf0>}
# Change the resolution of the cone and render again
>>> FindSource("Cone1").Resolution = 32
```



```
>>> Render()
```

You can also save state.

```
>>> from paraview.simple import *
>>> sph = Sphere()
>>> Render()
>>> servermanager.SaveState("/Users/berk/pythonstate.pvsm")
```

## Dealing with Time

If a reader or a filter supports time, it is easy to request a certain time step from Python. All time requests are set on views, which then propagate them to the representations which then propagate them to the visualization pipeline. Here is an example demonstrating how a time request can be made.

```
>>> Show(ExodusIIReader(FileName="../../../can.ex2"))
>>> Render()
# Get a nice view angle
>>> cam = GetActiveCamera()
>>> cam.Elevation(45)
>>> Render()
# Check the current view time
>>> view = GetActiveView()
>>> view.ViewTime
0.0
>>> reader = GetActiveSource()
>>> reader.TimestepValues
[0.0, 0.00010007373930420727, 0.00019990510190837085,
0.00029996439116075635, 0.00040008654468692839,
...]
>>> tsteps = reader.TimestepValues
# Let's be fancy and use a time annotation filter. This will show the
# current time value of the reader as text in the corner of the view.
>>> annTime = AnnotateTimeFilter(reader)
# Show the filter
>>> Show(annTime)
# Look at a few time steps. Note that the time value is requested not
# the time step index.
>>> view.ViewTime = tsteps[2]
>>> Render()
>>> view.ViewTime = tsteps[4]
>>> Render()
```

## Animating

Server Manager has a complicated animation engine based on keyframes and scenes. This section will introduce a few simple ways of animating your visualization. If you have a time-aware reader, you can animate it with `AnimateReader()`.

```
>>> reader = ExodusIIReader(FileName="../../../can.ex2")
>>> Show(reader)
>>> Render()
>>> c = GetActiveCamera()
>>> c.Elevation(95)
# Animate over all time steps. Note that we are not passing the optional
# 3rd argument here. If you pass a filename as the 3rd argument,
# AnimateReader will create a movie.
>>> AnimateReader(reader)
# Save the animation to an avi file
>>> AnimateReader(reader, filename="../../../movie.avi")
```

To animate properties other than time, you can use regular keyframes.

### **ParaView 3.8.0 and earlier**

Although the following script will work with 3.8.1 and later, it's not the recommended way since the changes done so will not be reflected in the GUI. Refer to the following sub-section for the recommended style for 3.8.1 and later versions.

```
>>> Sphere()
>>> Show()
>>> Render()

# Create an animation scene
>>> scene = servermanager.animation.AnimationScene()
# Add one view
>>> scene.ViewModules = [GetActiveView()]

# Create a cue to animate the StartTheta property
>>> cue = servermanager.animation.KeyFrameAnimationCue()
>>> cue.AnimatedProxy = GetActiveSource()
>>> cue.AnimatedPropertyName = "StartTheta"
# Add it to the scene's cues
>>> scene.Cues = [cue]

# Create 2 keyframes for the StartTheta track
>>> keyf0 = servermanager.animation.CompositeKeyFrame()
>>> keyf0.Interpolation = 'Ramp'
# At time = 0, value = 0
>>> keyf0.KeyTime = 0
>>> keyf0.KeyValues = [0]

>>> keyf1 = servermanager.animation.CompositeKeyFrame()
# At time = 1.0, value = 200
```

```
>>> keyf1.KeyTime = 1.0
>>> keyf1.KeyValues= [200]

# Add keyframes.
>>> cue.KeyFrames = [keyf0, keyf1]

>>> scene.Play()

# Some properties you can change
#
# Number of frames used in Sequence mode
# scene.NumberOfFrames = 100
#
# Or you can use real time mode
# scene.PlayMode = 'Real Time'
# scene.Duration = 20
```

---

**ParaView 3.8.1 onwards**

The following script will only work with ParaView versions 3.8.1 and later. It is now the recommended way for accessing animation scene and animation tracks since the updates are reflected in the GUI when running through the python shell from the ParaView application is used.

```
>>> Sphere()
>>> Show()
>>> Render()

# Get the application-wide animation scene
>>> scene = GetAnimationScene()

# Get the animation track for animating "StartTheta" on the active
source.
# GetAnimationTrack() creates a new track if none exists.
>>> cue = GetAnimationTrack("StartTheta")

# Create 2 keyframes for the StartTheta track
>>> keyf0 = CompositeKeyFrame()
>>> keyf0.Interpolation = 'Ramp'
# At time = 0, value = 0
>>> keyf0.KeyTime = 0
>>> keyf0.KeyValues= [0]

>>> keyf1 = CompositeKeyFrame()
# At time = 1.0, value = 200
>>> keyf1.KeyTime = 1.0
>>> keyf1.KeyValues= [200]

# Add keyframes.
>>> cue.KeyFrames = [keyf0, keyf1]
```

```
>>> scene.Play()

# Some properties you can change
#
# Number of frames used in Sequence mode
# scene.NumberOfFrames = 100
#
# Or you can use real time mode
# scene.PlayMode = 'Real Time'
# scene.Duration = 20
```

## GetAnimationTrack Usages

```
# Typical usage
>>> track = GetAnimationTrack("Center", 0, sphere) or
>>> track = GetAnimationTrack(sphere.GetProperty("Radius")) or

# this returns the track to animate visibility of the active source in
# all views.
>>> track = GetAnimationTrack("Visibility")

# For animating properties on implicit planes etc., use the following
# signatures:
>>> track = GetAnimationTrack(slice.SliceType.GetProperty("Origin"), 0) or
>>> track = GetAnimationTrack("Origin", 0, slice.SliceType)
```

## Loading Data Files

As seen throughout this document, you can always load a data file by explicitly creating the reader that can read the data file as follows:

```
>>> reader = ExodusIIReader(FileName="../../../can.ex2")
```

Alternatively, starting with ParaView 3.8, you can use `OpenDataFile()` function to let ParaView pick a reader using the extension of the file.

```
>>> reader = OpenDataFile("../../can.ex2")
```

## Writing Data Files (ParaView 3.9 or later)

To create a writer to write the output from a source, one can use the following:

```
from paraview.simple import *

# Specifying the source explicitly
>>> writer= CreateWriter("/.../filename.vtk", source)

# Using the active source
>>> writer= CreateWriter("/.../filename.vtk")
```

```
# Writing data from a particular output port
>>> writer= CreateWriter("../filename.vtk",
servermanager.OutputPort(source, 1))

# Now one change change the ivars on the writer

# To do the actual writing, use:
>>> writer.UpdatePipeline()
```

## Exporting CSV Data

To export a csv from the cell or point data associated with a source, one can use the following:

```
>>> writer = CreateWriter("../foo.csv", source)
>>> writer.FieldAssociation = "Points" # or "Cells"
>>> writer.UpdatePipeline()
>>> del writer
```

---

ParaView: [Welcome | Site Map <sup>[7]</sup>]

---

# Tools for Python Scripting

---

## New tools with python

If ParaView has been compiled with the Python wrapping, some advanced features become available to the user such as:

- Accessing a Python Shell
- Trace recording
- Macros
- Save ParaView state as a Python script

Those features can be reached from the **Tools** menu for the Shell and Trace access.

The management of macros is done inside the **Macros** menu.

To save the state as a Python script, just go to the **File** menu and choose **Save state** without forgetting to switch the file type to be **Python \*.py**.

## Macros

Macros allow the user to define a Python script as built-in actions that become accessible from the **Macros** menu or directly inside the Toolbar.

Once a local file is defined as a Macro (by clicking on **Create New Macro**) the given file is copied inside the user specific ParaView directory. No reference is kept to the original file. Therefore, if you keep editing the original file, the changes won't affect the macro itself.

The macros list is build dynamically at the ParaView startup by listing the content of the ParaView/Macros directory as well as the user specific ParaView/Macros directory. Note: if you want to rename a macro just rename the file in one of the given directory.

---

Deleted Macro **DO NOT** delete the file, but just renamed with a "." before the original filename and the file stay in the same directory.

Macros are displayed in the macros menu and the macros toolbar. The macros toolbar can be shown/hidden from the paraview main menu: View --> Toolbars --> Macro Toolbar. The toolbar may be hidden by default.

Note: python is not initialized until you open the python shell for the first time. If you run a macro from the macro toolbar or menu before the python shell has been opened for the first time, you will notice a slight delay as python initializes itself. You should see a wait cursor while python is initializing.

**CAUTION:** If this section is not relevant for your ParaView version, please read the previous version of that page in history. This is for ParaView 3.10

## Trace

Trace as been introduced in ParaView 3.6.2 as a new module. It can be imported with 'from paraview import smtrace', but normally the user never needs to directly use the trace module. The trace menu items provides everything for controlling trace:

- Start trace - starts trace. If an active trace was previously started, it will be stopped, cleared, and restarted.
- Stop trace- stops trace. The trace output generated so far will not be cleared until trace is started again or the python shell is reset. Some internal trace data structures hold references to c++ objects, so if you want to make sure everything is cleaned up try resetting the shell.
- Edit trace- opens the built in editor, creates a new document and fills it with the current trace output
- Save trace- opens a prompt for the user to specify a filename and saves trace to disk.

**TIP:** It's a good idea to stop trace before executing a trace script you just recorded. What I do is click the Disconnect server button in the paraview toolbar. This will clear the current pipeline objects, stop trace, and reset the python interpreter.

**TIP:** Trace only records property values as they are modified. If you have initialized a render view or a color lookup table prior to starting trace then the generated trace script may not perfectly reflect the state of the view or lookup table. For best results, start trace before performing any other actions. Also, see the CaptureAllProperties flag in the Trace Verbosity section below.

## Trace verbosity

Because of the way the GUI initializes certain proxies, some parts of the trace will be more verbose than others. For example, every time a data representation is created, seven or eight properties related to selection are modified. It might be a nice feature to provide the user with a way to specify property suppressions.

Trace can run with different verbose levels. In full property verbose mode, trace records all properties of a proxy when the proxy is registered. Normally, trace only records properties when they are modified from their default values. Control over the verbosity level is not currently available in the gui, but you can start trace manually in verbose mode using the python shell:

```
from paraview import smtrace
smtrace.start_trace(CaptureAllProperties=True)

# actions in the gui

smtrace.stop_trace()
```

## Known issues with trace

It is possible to perform actions in the GUI that do not translate properly to trace. Here is a list of things that do not currently work:

- views other than 3D render view (also have not tested mpi render view)
- time
- the start/stop trace buttons are not synced if trace is started manually from python

## New built in editor

For convenience there is a new built in script editor. When the editor is opened on OSX, the editor's menubar temporarily replaces paraview's menubar. If this becomes too obtrusive the menu could be replaced by toolbar buttons. It might be a nice feature to allow the user to specify a command to launch an external editor.

## New C++ API

New classes have been introduced under Qt/Python. They depend on classes in Qt/Core but not Qt/Components. A new class named pqPythonManager is available globally through the pqApplicationCore instance:

```
pqPythonManager* manager = qobject_cast<pqPythonManager*>(
    pqApplicationCore::instance()->manager("PYTHON_MANAGER"));
```

The python manager has a public method:

```
pqPythonDialog* pythonShellDialog();
```

Calling this method will return the python shell. Python will be initialized on the first call to this method and the returned pqPythonDialog will be ready to use. pqPythonDialog offers public methods for executing python files and strings of python code.

If you plan to call methods in the python c api, you must make the python interpreter active:

```
pqPythonDialog* dialog = manager->pythonShellDialog();
dialog->shell()->makeCurrent();

// Calls to python c api

dialog->shell()->releaseControl();
```

When the python interpreter is reset, pqPythonDialog emits a signal 'interpreterInitialized()'. The pqPythonManager listens for this signals and imports the paraview modules. So when this signal is triggered, it is not guaranteed that paraview python modules have been imported yet. After the paraview python modules are imported, the pqPythonManager emits a signal 'paraviewPythonModulesImported()'.

# Batch Processing

---

## Batch Processing

ParaView's **pvbatch** and **pvpython** command line executables substitute a python interpreter for the Qt GUI interface that most users control ParaView's back end data processing and rendering engine through. Either may be used for batch processing, that is to replay Visualization sessions in an exact, easily repeated way. The input to either comes in the form of the same python script that was described in the previous section.

Of the two, **pvbatch** is more specialized for batch processing and suited to running in an offline mode on dedicated data processing supercomputers because:

- It does not take in commands from the terminal, which is usually unavailable on this class of machines.

Therefore you must supply a filename of the script you want pvbatch to execute.

- It is permanently joined to the backend server and thus does not require TCP socket connections to it.

Therefore in the scripts that you give to pvbatch it is not possible to **Disconnect()** from the paired server or **Connect()** to a different one.

- It can be run directly as an MPI parallel program in which all pvbatch processes divide up the work and cooperate.

Therefore you typically start pvbatch like this:

```
[mpiexec -N <numprocessors>] pvbatch [args-for-pvbatch] script-filename [args-for-script]
```

## Creating the Input Deck

There are at least three ways to create a batch script.

The hardest one is writing it by hand using the syntax described in the previous section. You can of course use any text editor for this but you will probably be more productive if you set up a more fully featured python IDE like Idle or the python shell within the ParaView GUI so that you have access to interactive documentation, tab completion and quick preview capabilities. Another alternative is to let the ParaView GUI client record all of your actions into a python script by using the **Python Trace** feature. Later you can easily tweak the recorded script once you become familiar with ParaView's python syntax. The third, and to longtime ParaView users the most traditional way, is to instead record a ParaView state file and then load that via a small python script as demonstrated in the first example below.

## Examples

### Loading a state file and saving a rendered result

```
>>> from paraview.simple import *
# Load the state
>>> servermanager.LoadState("/Users/berk/myteststate.pvsm")
```

At this point you have a working pipeline instantiated on the server which you can use introspection on to access and then arbitrarily control anything within. At the core ParaView's is a visualization engine so we will demonstrate by simply generate and saving an image.

```
# Make sure that the view in the state is the active one so we don't
have to refer to it by name.
```



```
>>> SetActiveView(GetRenderView())
# Now render and save.
>>> Render()
>>> WriteImage("/Users/berk/image.png")
```

## parameter study

Parameter studies are one example of how batch processing can be extremely useful. In a parameter study one or more pipeline parameters (a filename, a timestep, or a filter property for example) are varied across some range but an otherwise identical script is replayed numerous times and results are saved. After the suite of sessions complete the set of results are easy to compare. For this type of work I recommend writing a higher level script that varies the parameter and for each value spawns off a pvbatch session where the parameter gets passed in as an argument to the ParaView python script.

The following is a slightly condensed version of a hierarchical set of scripts written during a benchmark study. This benchmark is an example of a parameter study in which the number of triangles rendered in the scene is varied and afterward we examine the output to determine how the rendering rate differs as a function of that parameter change.

This top level script varies the number of triangles and then submits parallel jobs to the cluster's PBS batch queue. See the qsub manpages or ask your system administrators for the exact syntax of the submission command.

```
RUNID=0
NNODES=8
TLIMIT=10
for NUMTRIS in 10 20 30 40 50
do
    mkdir ~/tmp/run${RUNID}

    qsub -N run${RUNID} \
        -l "walltime=0:${TLIMIT}:0.0" \
        select=${NNODES}:ncpus=8:arch=wds024c" \
        -j eo -e ~/tmp/run${ID}/outstreams.log \
        -v "RUNID=${ID} NNODES=${NNODES} NUMTRIS=${NUMTRIS}" \
        ~/level2.sh

    let RUNID+=1
done
```

The second level script is executed whenever it gets to the top of PBS's priority queue. It examines the parameters it is given and then runs paraview's pvbatch executable with them. It also does some bookkeeping tasks that are helpful when debugging the batch submission process.

```
echo "RUN NUMBER=${RUNID}"

#setup MPI environment
source ${HOME}/openmpipaths.sh

#prepare and run the parallel pvbatch program for the parameter value
we are given
batch_command="${HOME}/ParaView-3.8.1/build/bin/pvbatch
${HOME}/level3.py -# ${RUNID} -nt ${NUMTRIS}"
```

```
mpirun -np $NNODES --hostfile $PBS_NODEFILE $batch_command

#move the results to more permanent storage
mv /tmp/bench* ${HOME}/tmp/run${DDM_RUNNUM}
```

The final level is the script that is executed by pvbatch.

```
from paraview.simple import *
from optparse import OptionParser
import paraview.benchmark
import math
import sys
import time

parser = OptionParser()
parser.add_option("-#", "--runid", action="store",
dest="runid", type="int",
                default=-1, help="an identifier for this run")
parser.add_option("-nt", "--triangles", action="store",
dest="triangles", type="int",
                default=1, help="millions of triangles to render")
(options, args) = parser.parse_args()

print "#####"
print "RUNID = ", options.runid
print "START_TIME = ", time.localtime()
print "ARGS = ", sys.argv
print "OPTIONS = ", options
print "#####"

paraview.benchmark.maximize_logs()

TS = Sphere()
side=math.sqrt(options.triangles*1000000/2)
TS.PhiResolution = side
TS.ThetaResolution = side

dr = Show()
view.UseImmediateMode = 0
view = Render()

cam = GetActiveCamera()
for i in range(0,50):
    cam.Azimuth(3)
    Render()
    WriteImage('/tmp/bench_%d_image_%d.jpg' % (options.runid, i))

print "total Polygons:" +
```

```

str(dr.SMPProxy.GetRepresentedDataInformation(0).GetPolygonCount())
print "view.ViewSize:" + str(view.ViewSize)

paraview.benchmark.get_logs()
logname="/tmp/bench_" + str(options.runid) + "_rawlog.txt"
paraview.benchmark.dump_logs(logname)

print "#####"
print "END_TIME = ", time.localtime()

```

## large data example

Another important example is for visualizing extremely large datasets that can not be easily worked with interactively. In this setting, the user first constructs a visualization of a small but representative data set. Typically this takes place by recording a session in the standard GUI client running on some small and easily accessed machine. Later, the user changed the filename property of the reader in the recorded session file. Finally the user submits the script to a larger machine which performs the visualization offline and saves results for later inspection.

The essential thing that you need to be able to do for this is to substitute the filename and location of the original small dataset with the name and locations of the large one. There are two ways to do this.

The first way is to directly edit the filename in either the ParaView state file or the python script where it is loaded. The task is made easier by the fact that all readers conventionally name the input file name property "FileName". Standard python scripts are well described in other sections so we will describe paraview state files here instead. A paraview state file has the extension **.pvsm** and the internal format is a text based XML file. Simply open the pvsm file in a text editor, search for **FileName** and replace all occurrences of the old with the new.

For reference, the portion of a pvsm file that specifies a reader's input file is:

```

<Proxy group="sources" type="LegacyVTKFileReader" id="160" servers="1">
  <Property name="FileNameInfo" id="160.FileNameInfo" number_of_elements="1">
    <Element index="0" value="/Data/molar.vtk"/>
  </Property>
  <Property name="FileNames" id="160.FileNames" number_of_elements="1">
    <Element index="0" value="/Data/molar.vtk"/>
    <Domain name="files" id="160.FileNames.files"/>
  </Property>
  <Property name="TimestepValues" id="160.TimestepValues"/>
  <SubProxy name="Reader" servers="1"/>
</Proxy>

```

The second way is to set up the pipeline and then use introspection to find and then change the filename. This approach is easier to parameterize but somewhat more fragile since not all readers respond well to having their names changed once established. You should at least use caution and try to change the filename before the pipeline first runs. Otherwise more readers will be confused and you will also waste time processing the smaller file. When loading state files the proper place to do this is immediately after the **LoadState** command. For python scripts the place to do this is as near to the creation of the reader as possible, and certainly before any **Update** or **Render** commands. An example of how to do this follows:

```

>>> from paraview.simple import *
# Load the state
>>> servermanager.LoadState("/Users/berk/myteststate.pvsm")

```

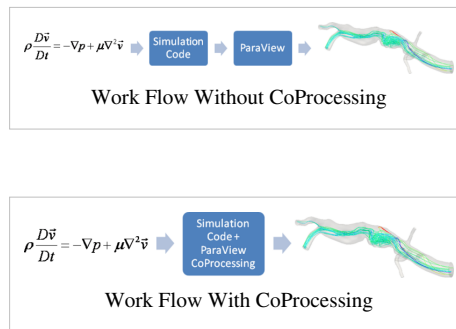
```
# Now the pipeline will be instantiated but it will not have updated
yet.
# You can programmatically obtain the reader from the pipeline starting
with this command, which lists all readers, sources and filters in the
pipeline.
>>> GetSources()
#{('box.ex2', '274'): <paraview.servermanager.ExodusIIReader object at 0x21b3eb70>}
# But it is easier if you note that readers are typically named
according to the name of the file that they are created for.
>>> reader = FindSource('box.ex2')
#Now you can change the filename with these two commands:
>>> reader.FileName = ['/path_to/can.ex2']
>>> reader.FileNameChanged()
```

# In-Situ/CoProcessing

## CoProcessing

### Background

Several factors are driving the growth of simulations. Computational power of computer clusters is growing while the price of individual computers is decreasing. Distributed computing techniques allow hundreds or even thousands of computer nodes to participate in a single simulation. The benefit of this computational power is that simulations are getting more accurate and useful for predicting complex phenomena. The downside to this growth in computational power is that enormous amounts of data need to be saved and analyzed to determine the results of the simulation. The ability to generate data has outpaced our ability to save and analyze the data. This bottle neck is throttling our ability to benefit from our improved computing resources. Simulations save their states only very infrequently to minimize storage requirements. This coarse temporal sampling makes it difficult to notice some complex behavior. To get past this barrier, ParaView can now be easily used to integrate post-processing/visualization directly with simulation codes. This feature is called co-processing in ParaView and the difference between work flows when using co-processing can be seen in the differences below.



### Technical Objectives

The main objective of the co-processing tool set is to integrate core data processing with the simulation to enable scalable data analysis while also being simple to use for the analyst. The tool set has two main parts:

- **An extensible and flexible co-processing library.** The co-processing library was designed to be flexible enough to be embedded in various simulation codes with relative ease. This flexibility is critical, as a library that requires a lot of effort to embed cannot be successfully deployed in a large number of simulations. The co-processing library also is easily extended so that users can easily deploy new analysis and visualization techniques to existing co-processing installations.
- **Configuration tools for co-processing configuration.** It is important for the users to be able to configure the co-processor using graphical user interfaces that are part of their daily work-flow.

Note that all of this must be done for large data. The co-processing library will almost always be run on a distributed system. For the largest simulations the visualization of extracts may also require a distributed system (i.e. a visualization cluster).

## Build Directions

As mentioned above, the two components for doing co-processing are a client side plugin and a server side library. It is recommended that they be compiled separately but from the same ParaView source revision/release. This is because currently the client side configuration tools outputs python code for the server side library to use. Differing ParaView versions may have different python interfaces to objects causing simulation co-processed runs to fail.

### ParaView Coprocessing Script Generator Plugin

The plugin for generating python scripts for co-processing is a client side plugin. The CMake option to turn on the script generator plugin is `PARAVIEW_BUILD_PLUGIN_CoProcessingScriptGenerator`. Note that since this is a client side plugin that the `PARAVIEW_BUILT_QT_GUI` option must be on.

### CoProcessing Library

The directions for building the co-processing library can be a bit more complex. We assume that it will be built on a cluster or supercomputer. Complexities may arise from having to build mesa, use off-screen rendering, build static libraries, and/or cross-compiling <sup>[1]</sup>. We won't go into those details here though but refer interested people to the ParaView <sup>[2]</sup> and VTK <sup>[3]</sup> wikis.

The additional flags that should be turned on for coprocessing include:

- **BUILD\_SHARED\_LIBS** Should normally be set to ON unless you're compiling on a machine that doesn't support shared libs (e.g. IBM BG/L).
- **PARAVIEW\_ENABLE\_PYTHON** Set to ON in order to use the scripts created from the coprocessing plugin. Note that of the 3.10 release that this is not required to be on though.
- **PARAVIEW\_USE\_MPI** Set the ParaView server to use MPI to run in parallel. Also check that the proper version of MPI is getting used.
- **CMAKE\_BUILD\_TYPE** Should be set to *Release* in order to optimize the ParaView code.
- **PARAVIEW\_ENABLE\_COPROCESSING** The CMake option to build the coprocessing libraries.
- **BUILD\_COPROCESSING\_ADAPTORS** The CMake option to turn on code that may be useful for creating the simulation code adaptor.
- **BUILD\_FORTTRAN\_COPROCESSING\_ADAPTORS** The CMake option to turn on code that may be useful for creating an adaptor for a simulation code written in Fortran or C.
- **PYTHON\_ENABLE\_MODULE\_vtkCoProcessorPython** On by default.
- **BUILD\_PYTHON\_COPROCESSING\_ADAPTOR** Experimental code.

If ParaView was build with python enabled, the co-processing library can be tested on the installed system with the CTest command `ctest -R CoProcessing`. The `CoProcessingTestPythonScript` test does an actual coprocessing run and the `CoProcessingPythonScriptGridPlot` and `CoProcessingPythonScriptPressurePlot` test verify the output is correct. In addition, looking at the code for this test is a useful beginning point for learning to use the co-processing library.

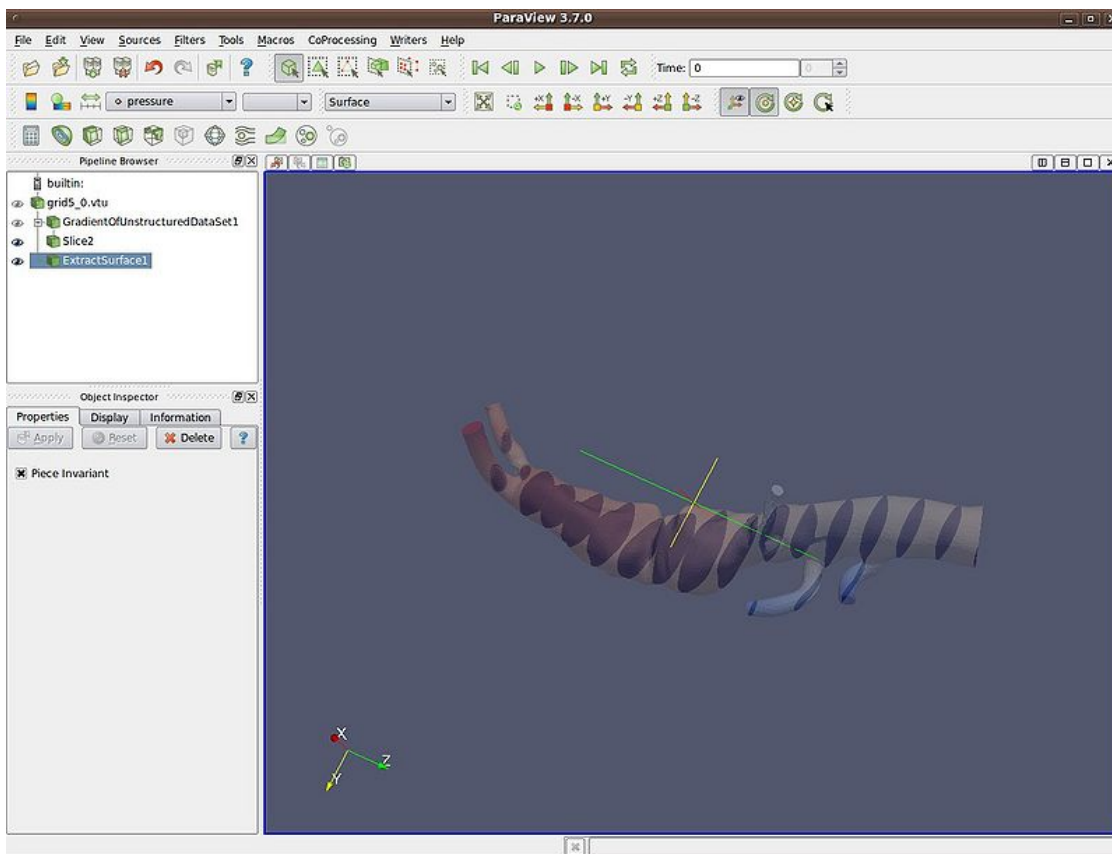
If you are using Mesa, the following options should be used:

- **VTK\_OPENGL\_HAS\_OSMESA** Set to ON.
- **OSMESA\_INCLUDE\_DIR**, **OPENGL\_xmesa\_INCLUDE\_DIR** and **OPENGL\_INCLUDE\_DIR** Make sure this is not set to the OpenGL version of the header files.
- **OSMESA\_LIBRARY**, **OPENGL\_gl\_LIBRARY** and **OPENGL\_glu\_LIBRARY** Set these to the proper libraries.
- **VTK\_USE\_OFFSCREEN** Set to ON

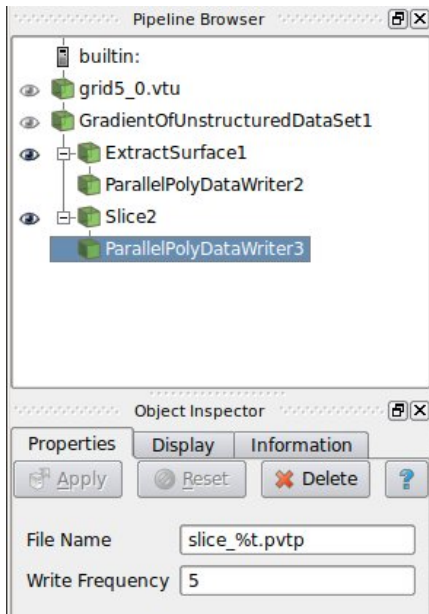
## Running the CoProcessing Script Generator

After starting the ParaView client, go to Tools->Manage Plugins, highlight CoProcessingPlugin and click Load Selected. Alternatively, if you expect to be using this plugin often you can tell ParaView to automatically load the plugin by clicking to the left of CoProcessingPlugin to expand the list and choose the Auto Load option. This will add a *CoProcessing* and *Writers* menu to the GUI.

Once the plugin is loaded, load a simulation result into ParaView that is similar to the run you wish to do co-processing with. Normally the result will be for a discretization of the same geometry and contain the same field variables (or at least all of the field variables you wish to examine) as the simulation grid that co-processing will be performed on. The only difference between these two grids will be that the grid used for the co-processing run will be at a much higher resolution. In the picture below, a pipeline is shown for a result from a PHASTA (A CFD code led by Ken Jansen at the University of Colorado) simulation of incompressible flow through an aortic aneurysm. The simulation results include the velocity and pressure fields and was done on a grid with 235,282 cells and 45,175 points. It is important to use the same names for the field arrays as will be used in the grids the coprocessing library will be working with. The filters specify which fields to use by their name and problems will occur if the filters cannot find the correct array. The pipeline computes the vorticity, takes 20 slices through the domain, and extracts the surface of the grid.



Once the pipeline is set up, the writers must be added to use for the filters that results are desired from. This is done by selecting the appropriate writer to output the results from each desired filter. In this example a *ParallelPolyDataWriter* is created for the Slice filter and the ExtractSurface filter. For the writer for the Slice filter we set the name of the file to be *slice\_%t.pvtp* where *%t* will get replaced with the time step each time it is output. We also set the write frequency to be every fifth time step. The picture below shows this.

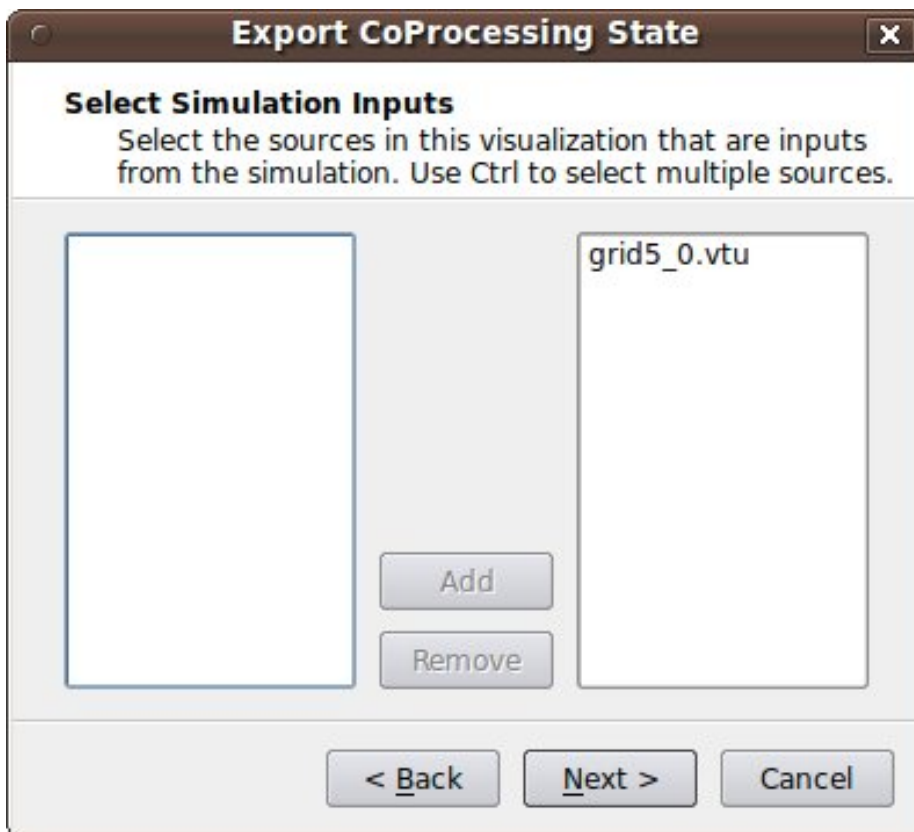


Similarly, we do the same for the writer for the ExtractSurface filter but we want to use a different file name and we can set a different write frequency if desired.

Note that when we hit enter we get the error message *Cannot show the data in the current view although the view reported that it can show the data. This message can be ignored and is already entered as a bug to fix in ParaView.*

The final step is to go through the CoProcessing->Export State wizard to associate the grid with a name that the python script uses and specify a file name for the script. The steps are:

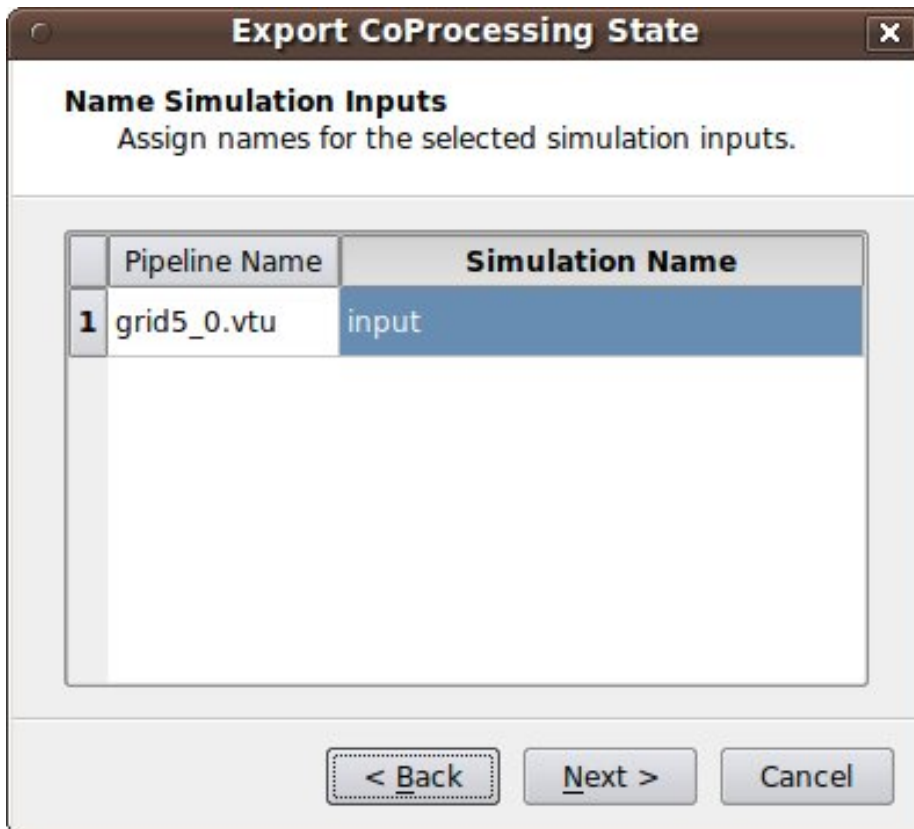
1. **Export Co-Processing State** Click Next to proceed.
2. **Select Simulation Inputs** Choose which grids to use by highlighting the grid and clicking Add, then click Next.



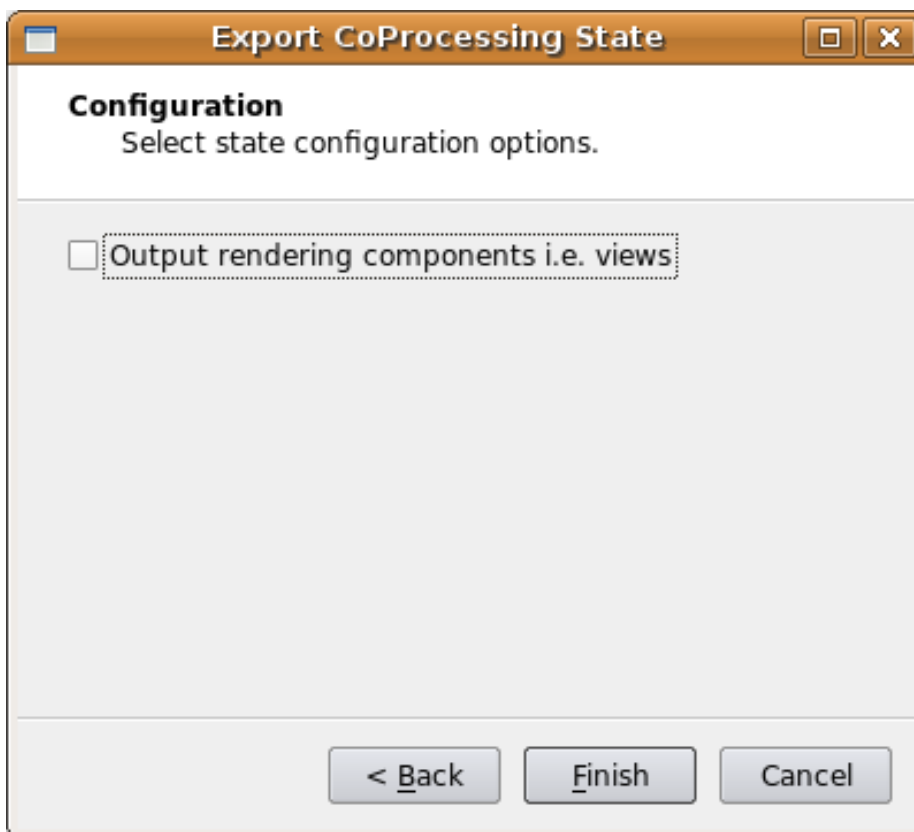
3. **Name Simulation Inputs** Click in the boxes under *Simulation Name* and change the name to what you have named the grid in your adaptor. Our convention is to use "input" for simulations with a single grid or multiblock data



set.

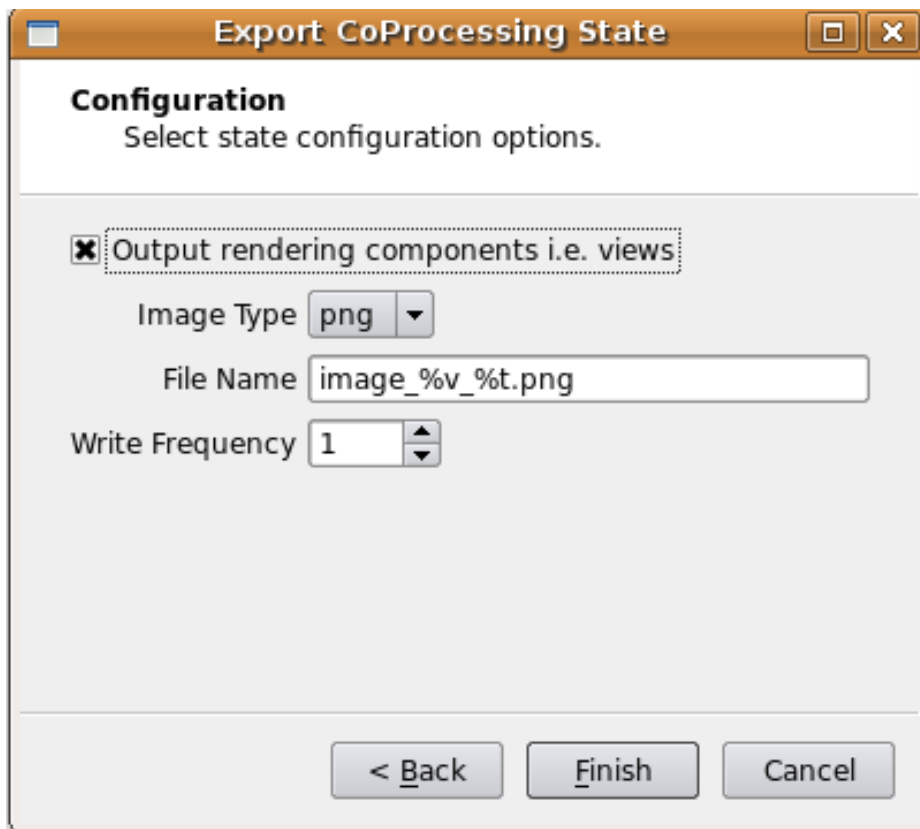


4. **Configuration** We usually want the full data and not just a picture of the results so for this we leave "Output rendering components i.e. views" unchecked and just hit Finish.



If we do want to output images in addition to any writers we may have specified, check "Output rendering components i.e. views" and fill in the proper information. Note that the image type is specified based on the file

name extension and that %t will be replaced by the actual time step in the output file name. Additionally, if there are multiple views then %v will be replaced by the view index in the output file name.



5. **Save Server State:** Specify the file name the script is to be saved as.

The resulting python script should look something like:

```
try: paraview.simple
except: from paraview.simple import *

cp_writers = []

def RequestDataDescription(datadescription):
    "Callback to populate the request for current timestep"
    timestep = datadescription.GetTimeStep()

    if (timestep % 5 == 0) or (timestep % 200 == 0) :

datadescription.GetInputDescriptionByName('input').AllFieldsOn()

datadescription.GetInputDescriptionByName('input').GenerateMeshOn()

def DoCoProcessing(datadescription):
    "Callback to do co-processing for current timestep"
    global cp_writers
    cp_writers = []
    timestep = datadescription.GetTimeStep()
```

```

grid5_0_vtu = CreateProducer( datadescription, "input" )

GradientOfUnstructuredDataSet1 = GradientOfUnstructuredDataSet(
guiName="GradientOfUnstructuredDataSet1", ScalarArray=['POINTS',
'veLOCITY'], ResultArray
Name='Vorticity', ComputeVorticity=1, FasterApproximation=0 )

Slice2 = Slice( guiName="Slice2",
SliceOffsetValues=[-10.23284210526316, -8.8684631578947375,
-7.5040842105263152, -6.1397052631578966, -4.77532631578947
43, -3.4109473684210529, -2.0465684210526316, -0.68218947368421201,
0.68218947368420935, 2.0465684210526316, 3.4109473684210521,
4.7753263157894743, 6.139705
263157893, 7.5040842105263152, 8.8684631578947375, 10.23284210526316,
11.597221052631578, 12.961600000000001], SliceType="Plane" )

SetActiveSource(GradientOfUnstructuredDataSet1)
ExtractSurface1 = ExtractSurface( guiName="ExtractSurface1",
PieceInvariant=1 )

SetActiveSource(Slice2)
ParallelPolyDataWriter3 = CreateWriter( XMLPPolyDataWriter,
"slice_%t.pvtp", 5 )

SetActiveSource(ExtractSurface1)
ParallelPolyDataWriter2 = CreateWriter( XMLPPolyDataWriter,
"surface_%t.pvtp", 200 )

Slice2.SliceType.Origin = [-2.8049160242080688, 2.1192346811294556,
1.7417440414428711]
Slice2.SliceType.Offset = 0.0
Slice2.SliceType.Normal = [0.0, 0.0, 1.0]

for writer in cp_writers:
    if timestep % writer.cpFrequency == 0:
        writer.FileName = writer.cpFileName.replace("%t",
str(timestep))
        writer.UpdatePipeline()

def CreateProducer(datadescription, gridname):
    "Creates a producer proxy for the grid"
    if not datadescription.GetInputDescriptionByName(gridname):
        raise RuntimeError, "Simulation input name '%s' does not exist" %
gridname

```

```

grid = datadescription.GetInputDescriptionByName(gridname).GetGrid()
producer = TrivialProducer()
producer.GetClientSideObject().SetOutput(grid)
producer.UpdatePipeline()
return producer

def CreateWriter(proxy_ctor, filename, freq):
    global cp_writers
    writer = proxy_ctor()
    writer.FileName = filename
    writer.add_attribute("cpFrequency", freq)
    writer.add_attribute("cpFileName", filename)
    cp_writers.append(writer)
    return writer

```

## Creating the Adaptor

The largest amount of work in getting co-processing working with a simulation code is usually creating the adaptor that can pass a VTK data set or composite data set with fields specified over it. A simple view of the program control is



By doing it like this, there is minimal modification to the simulation code (only calls to the adaptor like Initialize, Finalize, and CoProcess). The adaptor deals with the coprocessing library as well as creating VTK data structures from the simulation code as needed. The general flow for the simulation code and where it calls the adaptor would look like:

```

initialize coprocessor
for i in number of time steps
    compute simulation information at time step i
    call coprocessing library
        coprocessing library determines if any coprocessing needs to be performed at time step i
finalize coprocessor

```

The interactions between the adaptor and ParaView are mostly done through the `vtkCPPProcessor` class. Documentation for this and the other classes in the coprocessing library are in the ParaView online documentation <sup>[4]</sup> as well as grouped in a module <sup>[5]</sup>. Details on the specific steps are provided below.

## Initialize the Co-Processor

To initialize the co-processor, a `vtkCPPProcessor` object must be created. This object should not be deleted until the finalization step. After creating a `vtkCPPProcessor` object, you must call `vtkCPPProcessor::Initialize()` to initialize the coprocessing library. The next step is to create `vtkCPPPipeline` objects. These objects are used to represent the VTK pipelines that will be used to generate the output data. Most users will probably only use the `vtkCPPythonScriptPipeline` class that derives from `vtkCPPPipeline` since that pipeline is the one that is used with the

python scripts generated from the coprocessing plugin. `vtkCPPythonScriptPipeline::Initialize()` takes in the python script name with the full path and sets up the ParaView python interpreter.

## Call the Co-Processor to Execute the Desired Filters

Most users will have a single call to the adaptor at the end of each simulation time step. This will pass in the simulation data structures to create the VTK data structures from. The created VTK data structures will be an object that derives from `vtkDataObject`. First though we want to check if we actually need to do any co-processing during this call. The simulation code will not know this information. Because of this, the simulation code should call the adaptor every time step. The adaptor gets the information about what data is needed for executing the `vtkCPPipelines` for each call. This is done in `vtkCPProcessor::RequestDataDescription()` by going through all of the `vtkCPPipeline` objects and querying them for what is needed. `vtkCPProcessor::RequestDataDescription()` takes in a `vtkCPDataDescription` object that contains information about the current simulation time and time step. The method returns 0 if no coprocessing information needs to be computed during the call and 1 otherwise. If no coprocessing needs to be performed then the adaptor should return control back to the simulation code.

If co-processing is requested then the adaptor needs to make sure that the VTK grids and fields are up to date. Upon return from `vtkCPProcessor::RequestDataDescription()`, the `vtkCPDataDescription` object will have information on what is needed by the pipelines in order to successfully execute them during this call to the adaptor. This information is stored for each co-processing input (the data objects that are given identifier names in the plugin) and accessed through the `vtkCPDataDescription` as `vtkCPIInputDataDescription` objects. The `vtkCPIInputDataDescription` objects specify whether or not they require a grid and if so which fields they require. These objects are also used to pass the VTK data objects created in the adaptor to the pipeline by using the `vtkCPIInputDataDescription::SetGrid()` method. The final step in this call to the adaptor is then to perform the actual co-processing by calling the `vtkCPProcessor::CoProcess()` method.

## Finalize the Co-Processor

Finalization entails calling `vtkCPProcessor::Finalize()` and cleaning up any remaining VTK objects by calling `Delete()`;

---

While initially there may seem to be a significant hurdle to overcome in order to implement the adaptor for co-processing, the user only needs to know enough information to create the VTK data objects from their own simulation data structures. The adaptor can be made computationally efficient by taking advantage of intimate knowledge of the simulation code and knowing that VTK filters will not modify input data. Examples include:

- Only creating the VTK grid in the first call to the adaptor if the grid is not changing topologically or geometrically.
- Using the `SetArray()` method to build `vtkDataArray` objects that use existing memory used by the simulation code instead of allocating and copying existing memory.

Finally, we suggest looking at either the C++ coprocessing example and/or the python coprocessing example to see how the entire co-processing flow works. Additionally, the code in `ParaView3/CoProcessing/CoProcessor/Testing/Cxx/PythonScriptCoProcessingExample.cxx` is also worth a look. This is used to test the co-processing library and can be run with the command "ctest -R CoProcessing".

---

## References

- [1] [http://paraview.org/Wiki/Cross\\_compiling\\_ParaView3\\_and\\_VTK](http://paraview.org/Wiki/Cross_compiling_ParaView3_and_VTK)
- [2] <http://paraview.org/Wiki/ParaView>
- [3] <http://www.vtk.org/Wiki/VTK>
- [4] <http://www.paraview.org/ParaView3/Doc/Nightly/html/classes.html>
- [5] [http://www.paraview.org/ParaView3/Doc/Nightly/html/group\\_\\_CoProcessing.html](http://www.paraview.org/ParaView3/Doc/Nightly/html/group__CoProcessing.html)

## C++ CoProcessing example

---

This example is used to demonstrate how the co-processing library can be used with a C++ based simulation code. In the ParaView/CoProcessing/Adaptors/FortranAdaptors directory there is code useful for integrating C or Fortran based simulation codes with the co-processing library. Note that this example requires MPI to be available on your system. The executable takes in a python coprocessing script and a number of time steps to be run for. Note to remember to set your system environment properly. See [[1]] for details.

### CoProcessingExample.cxx

```
#include "vtkCPDataDescription.h"
#include "vtkCPInputDataDescription.h"
#include "vtkCPPProcessor.h"
#include "vtkCPPythonScriptPipeline.h"
#include "vtkElevationFilter.h"
#include "vtkMultiProcessController.h"
#include "vtkPolyData.h"
#include "vtkSmartPointer.h"
#include "vtkSphereSource.h"
#include "vtkXMLUnstructuredGridReader.h"

#include <mpi.h>
#include <string>

class DataGenerator {
public:
    DataGenerator()
    {
        this->Sphere = vtkSmartPointer<vtkSphereSource>::New();
        this->Sphere->SetThetaResolution(30);
        this->Sphere->SetPhiResolution(30);
        int procId;
        MPI_Comm_rank(MPI_COMM_WORLD, &procId);
        this->Sphere->SetCenter(procId*4.0, 0, 0);
        this->Elevation = vtkSmartPointer<vtkElevationFilter>::New();
        this->Elevation->SetInputConnection(this->Sphere->GetOutputPort());
        this->Index = 0;
    }

    vtkSmartPointer<vtkPolyData> GetNext()
```

```

{
    double radius = fabs(sin(0.1 * this->Index));
    this->Index++;
    this->Sphere->SetRadius(1.0 + radius);
    this->Elevation->Update();
    vtkSmartPointer<vtkPolyData> ret = vtkSmartPointer<vtkPolyData>::New();
    ret->DeepCopy(this->Elevation->GetOutput());
    return ret;
}

protected:
    int Index;
    vtkSmartPointer<vtkSphereSource> Sphere;
    vtkSmartPointer<vtkElevationFilter> Elevation;
};

int main(int argc, char* argv[])
{
    if (argc < 3)
    {
        printf("Usage: %s <cp python file> <number of steps>\n", argv[0]);
        return 1;
    }
    // we assume that this is done in parallel
    MPI_Init(&argc, &argv);

    std::string cpPythonFile = argv[1];
    int nSteps = atoi(argv[2]);

    vtkCPPProcessor* processor = vtkCPPProcessor::New();
    processor->Initialize();
    vtkCPPythonScriptPipeline* pipeline =
    vtkCPPythonScriptPipeline::New();

    // read the coprocessing python file
    if(pipeline->Initialize(cpPythonFile.c_str()) == 0)
    {
        cout << "Problem reading the python script.\n";
        return 1;
    }

    processor->AddPipeline(pipeline);
    pipeline->Delete();

    if (nSteps == 0)
    {
        return 0;
    }

```

```
}

// create a data source, typically this will come from the adaptor
// but here we use generator to create it ourselves
DataGenerator generator;

// do coprocessing
double tStart = 0.0;
double tEnd = 1.0;
double stepSize = (tEnd - tStart)/nSteps;

vtkCPDataDescription* dataDesc = vtkCPDataDescription::New();
dataDesc->AddInput("input");

for (int i = 0; i < nSteps; ++i)
{
    double currentTime = tStart + stepSize*i;
    // set the current time and time step
    dataDesc->SetTimeData(currentTime, i);

    // check if the script says we should do coprocessing now
    if(processor->RequestDataDescription(dataDesc) != 0)
    {
        // we are going to do coprocessing so use generator to
        // create our grid at this timestep and provide it to
        // the coprocessing library
        vtkSmartPointer<vtkDataObject> dataObject =
            generator.GetNext();

        dataDesc->GetInputDescriptionByName("input")->SetGrid(dataObject);
        processor->CoProcess(dataDesc);
    }
}

dataDesc->Delete();
processor->Finalize();
processor->Delete();

MPI_Finalize();

return 0;
}
```



## CMakeLists.txt

```
cmake_minimum_required(VERSION 2.6)

PROJECT(CoProcessingExample)

FIND_PACKAGE(ParaView REQUIRED)
INCLUDE(${PARAVIEW_USE_FILE})

FIND_PACKAGE(MPI REQUIRED)
INCLUDE_DIRECTORIES(${MPI_INCLUDE_PATH})

ADD_EXECUTABLE(CoProcessingExample CoProcessingExample.cxx)
TARGET_LINK_LIBRARIES(CoProcessingExample vtkCoProcessor)
```

## Python Scripts

The first python script below is used to just output the actual results of the example. This would correspond to a simulation run with a coarse grid in order to set up coprocessing runs for larger grids where outputting the entire simulation results can be computationally prohibitive.

```
try: paraview.simple
except: from paraview.simple import *

cp_writers = []

def RequestDataDescription(datadescription):
    "Callback to populate the request for current timestep"
    timestep = datadescription.GetTimeStep()
    input_name = 'input'
    if (timestep % 1 == 0) :

datadescription.GetInputDescriptionByName(input_name).AllFieldsOn()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOn()
    else:

datadescription.GetInputDescriptionByName(input_name).AllFieldsOff()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOff()

def DoCoProcessing(datadescription):
    "Callback to do co-processing for current timestep"
    global cp_writers
    cp_writers = []
    timestep = datadescription.GetTimeStep()

    Sphere1 = CreateProducer( datadescription, "input" )
    ParallelPolyDataWriter1 = CreateWriter( XMLPPolyDataWriter,
"input_grid_%t.pvtp", 1 )
```

```

    for writer in cp_writers:
        if timestep % writer.cpFrequency == 0:
            writer.FileName = writer.cpFileName.replace("%t",
str(timestep))
            writer.UpdatePipeline()

    # explicitly delete the proxies -- we do it this way to avoid
problems with prototypes
    tobedeleted = GetProxiesToDelete()
    while len(tobedeleted) > 0:
        Delete(tobedeleted[0])
        tobedeleted = GetProxiesToDelete()

def GetProxiesToDelete():
    iter = servermanager.vtkSMPProxyIterator()
    iter.Begin()
    tobedeleted = []
    while not iter.IsAtEnd():
        if iter.GetGroup().find("prototypes") != -1:
            iter.Next()
            continue
        proxy = servermanager._getPyProxy(iter.GetProxy())
        proxygroup = iter.GetGroup()
        iter.Next()
        if proxygroup != 'timekeeper' and proxy != None and
proxygroup.find("pq_helper_proxies") == -1 :
            tobedeleted.append(proxy)

    return tobedeleted

def CreateProducer(datadescription, gridname):
    "Creates a producer proxy for the grid"
    if not datadescription.GetInputDescriptionByName(gridname):
        raise RuntimeError, "Simulation input name '%s' does not exist" %
gridname
    grid = datadescription.GetInputDescriptionByName(gridname).GetGrid()
    producer = TrivialProducer()
    producer.GetClientSideObject().SetOutput(grid)
    producer.UpdatePipeline()
    return producer

def CreateWriter(proxy_ctor, filename, freq):
    global cp_writers
    writer = proxy_ctor()
    writer.FileName = filename
    writer.add_attribute("cpFrequency", freq)

```

```

writer.add_attribute("cpFileName", filename)
cp_writers.append(writer)
return writer

```

This second script is still rather simple and only performs a cut on the input from the simulation code. It demonstrates though how desired results can be obtained while performing coprocessing at specified time steps.

```

try: paraview.simple
except: from paraview.simple import *
cp_writers = []

def RequestDataDescription(datadescription):
    "Callback to populate the request for current timestep"
    timestep = datadescription.GetTimeStep()
    input_name = 'input'
    if (timestep % 5 == 0) :

datadescription.GetInputDescriptionByName(input_name).AllFieldsOn()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOn()
    else:

datadescription.GetInputDescriptionByName(input_name).AllFieldsOff()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOff()

def DoCoProcessing(datadescription):
    "Callback to do co-processing for current timestep"
    global cp_writers
    cp_writers = []
    timestep = datadescription.GetTimeStep()

    filename_0_pvtp = CreateProducer( datadescription, "input" )

    Clip2 = Clip( guiName="Clip2", InsideOut=0, UseValueAsOffset=0,
Scalars=['POINTS', 'Elevation'], Value=0.0, ClipType="Plane" )
    Clip2.ClipType.Normal = [0.0, 1.0, 0.0]
    Clip2.ClipType.Origin = [1.9999999105930328, 0.0, 0.0]
    Clip2.ClipType.Offset = 0.0

    ParallelUnstructuredGridWriter2 = CreateWriter(
XMLPUnstructuredGridWriter, "Cut_%.pvtu", 5 )
    for writer in cp_writers:
        if timestep % writer.cpFrequency == 0:
            writer.FileName = writer.cpFileName.replace("%t",
str(timestep))
            writer.UpdatePipeline()

```

```

    # explicitly delete the proxies -- we do it this way to avoid
    problems with prototypes
    tobdeleted = GetProxiesToDelete()
    while len(tobdeleted) > 0:
        Delete(tobdeleted[0])
        tobdeleted = GetProxiesToDelete()

def GetProxiesToDelete():
    iter = servermanager.vtkSMPProxyIterator()
    iter.Begin()
    tobdeleted = []
    while not iter.IsAtEnd():
        if iter.GetGroup().find("prototypes") != -1:
            iter.Next()
            continue
        proxy = servermanager._getPyProxy(iter.GetProxy())
        proxygroup = iter.GetGroup()
        iter.Next()
        if proxygroup != 'timekeeper' and proxy != None and
proxygroup.find("pq_helper_proxies") == -1 :
            tobdeleted.append(proxy)

    return tobdeleted

def CreateProducer(datadescription, gridname):
    "Creates a producer proxy for the grid"
    if not datadescription.GetInputDescriptionByName(gridname):
        raise RuntimeError, "Simulation input name '%s' does not exist" %
gridname
    grid = datadescription.GetInputDescriptionByName(gridname).GetGrid()
    producer = TrivialProducer()
    producer.GetClientSideObject().SetOutput(grid)
    producer.UpdatePipeline()
    return producer

def CreateWriter(proxy_ctor, filename, freq):
    global cp_writers
    writer = proxy_ctor()
    writer.FileName = filename
    writer.add_attribute("cpFrequency", freq)
    writer.add_attribute("cpFileName", filename)
    cp_writers.append(writer)
    return writer

```

## References

[1] [http://paraview.org/Wiki/ParaView/Python\\_Scripting#Getting\\_Started](http://paraview.org/Wiki/ParaView/Python_Scripting#Getting_Started)

# Python CoProcessing Example

---

This example is used to demonstrate how the co-processing library can be used with a python based simulation code. Note that this example requires MPI to be available on your system as well as pyMPI to initialize and finalize MPI from the python script. The executable takes in a python co-processing script and a number of time steps to be run for. Note to remember to set your system environment properly. See [[1]] for details.

## Python driver code

```
import sys
if len(sys.argv) != 3:
    print "command is 'pyMPI <python driver code> <script name> <number of time steps>'"
    sys.exit(1)
import mpi
mpi.initialized()
import paraview
import paraview.vtk as vtk

# initialize and read input parameters
paraview.options.batch = True
paraview.options.symmetric = True

def _refHolderMaker(obj):
    def _refHolder(obj2, string):
        tmp = obj
    return _refHolder

def coProcess(grid, time, step, scriptname):
    import vtkCoProcessorPython # import libvtkCoProcessorPython for
older PV versions
    if scriptname.endswith(".py"):
        scriptname = scriptname.rstrip(".py")
    try:
        cpscript = __import__(scriptname)
    except:
        print 'Cannot find ', scriptname, ' -- no coprocessing will be
performed.'
        return

    datadescription = vtkCoProcessorPython.vtkCPDataDescription()
    datadescription.SetTimeData(time, step)
    datadescription.AddInput("input")
    cpscript.RequestDataDescription(datadescription)
    inputdescription =
```

```

datadescription.GetInputDescriptionByName("input")
    if inputdescription.GetIfGridIsNecessary() == False:
        return

    inputdescription.SetGrid(grid)
    cpscript.DoCoProcessing(datadescription)

try:
    numsteps = int(sys.argv[2])
except ValueError:
    print 'the last argument should be a number'
    numsteps = 10

for step in range(numsteps):
    # assume simulation time starts at 0
    time = step/float(numsteps)

    # create the input to the coprocessing library. normally
    # this will come from the adaptor
    imageData = vtk.vtkImageData()
    imageData.SetOrigin(0, 0, 0)
    imageData.SetSpacing(.1, .1, .1)
    imageData.SetDimensions(10, 12, 12)
    pointArray = vtk.vtkDoubleArray()
    pointArray.SetNumberOfTuples(imageData.GetNumberOfPoints())
    for i in range(imageData.GetNumberOfPoints()):
        pointArray.SetValue(i, i)
    pointArray.SetName("pointData")
    imageData.GetPointData().AddArray(pointArray)

    # "perform" coprocessing. results are outputted only if
    # the passed in script says we should at time/step
    coProcess(imageData, time, step, sys.argv[1])

mpi.finalized()

```

## Sample coprocessing script

```

try: paraview.simple
except: from paraview.simple import *

cp_writers = []

def RequestDataDescription(datadescription):
    "Callback to populate the request for current timestep"
    timestep = datadescription.GetTimeStep()
    input_name = 'input'
    if (timestep % 1 == 0) :

```

```

datadescription.GetInputDescriptionByName(input_name).AllFieldsOn()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOn()
    else:

datadescription.GetInputDescriptionByName(input_name).AllFieldsOff()

datadescription.GetInputDescriptionByName(input_name).GenerateMeshOff()

def DoCoProcessing(datadescription):
    "Callback to do co-processing for current timestep"
    global cp_writers
    cp_writers = []
    timestep = datadescription.GetTimeStep()

    Sphere1 = CreateProducer( datadescription, "input" )
    ImageWriter1 = CreateWriter( XMLPImageDataWriter,
    "input_grid_%t.pvti", 1 )

    for writer in cp_writers:
        if timestep % writer.cpFrequency == 0:
            writer.FileName = writer.cpFileName.replace("%t",
str(timestep))
            writer.UpdatePipeline()

    # explicitly delete the proxies -- we do it this way to avoid
problems with prototypes
    tobedeleted = GetProxiesToDelete()
    while len(tobedeleted) > 0:
        Delete(tobedeleted[0])
        tobedeleted = GetProxiesToDelete()

def GetProxiesToDelete():
    iter = servermanager.vtkSMPProxyIterator()
    iter.Begin()
    tobedeleted = []
    while not iter.IsAtEnd():
        if iter.GetGroup().find("prototypes") != -1:
            iter.Next()
            continue
        proxy = servermanager._getPyProxy(iter.GetProxy())
        proxygroup = iter.GetGroup()
        iter.Next()
        if proxygroup != 'timekeeper' and proxy != None and
proxygroup.find("pq_helper_proxies") == -1 :
            tobedeleted.append(proxy)

```

```
        return tobedeleted

def CreateProducer(datadescription, gridname):
    "Creates a producer proxy for the grid"
    if not datadescription.GetInputDescriptionByName(gridname):
        raise RuntimeError, "Simulation input name '%s' does not exist" %
gridname
    grid = datadescription.GetInputDescriptionByName(gridname).GetGrid()
    producer = TrivialProducer()
    producer.GetClientSideObject().SetOutput(grid)
    producer.UpdatePipeline()
    return producer

def CreateWriter(proxy_ctor, filename, freq):
    global cp_writers
    writer = proxy_ctor()
    writer.FileName = filename
    writer.add_attribute("cpFrequency", freq)
    writer.add_attribute("cpFileName", filename)
    cp_writers.append(writer)
    return writer
```



---

# Plugins

---

## What are Plugins?

---

### Introduction

ParaView comes with plethora of functionality bundled in: several readers, multitude of filters, quite a few different types of views etc. However, it is not uncommon for developers to add new functionality to ParaView. For example to add support for their new file format, incorporate a new filter, etc. ParaView makes it possible to add new functionality by using an extensive plugin mechanism.

Plugins can be used to extend ParaView in several ways:

- Add new readers
- Add new writers
- Add new filters
- Add new GUI components
- Add new views
- Add new representations

### Plugin Types

Plugins are distributed as shared libraries (\*.so on Unix, \*.dylib on Mac, \*.dll on Windows etc). For a plugin to be loadable in ParaView, it must be built with the same version of ParaView as it is expected to be deployed on. Plugins can be classified into two broad categories:

- Server-side plugins

These are plugins that extend the algorithmic capabilities for ParaView eg. new filters, readers, writers etc. Since in ParaView data is processed on the server-side, these plugins need to be loaded on the server.

- Client-side plugins

These are plugins that extend the ParaView GUI eg. property panels for new filters, toolbars, views etc. These plugins need to be loaded on the client.

Oftentimes a plugin has both server-side as well as client-side components to it eg. a plugin that adds a new filter and a property panel that goes with that filter. Such plugins need to be loaded both on the server as well as the client.

Generally, users don't have to worry whether a plugin is a server-side or client-side plugin. Simply load the plugin on the server as well as the client. ParaView will include relevant components from plugin on each of the processes.

---

# Included Plugins

---

## Included Plugins

ParaView comes with a collection of plugins that the community has developed.

- Adios

Loads pixie format files written with the Adios library. This reader support the staging capability of Adios which allow the user to pipe simulation kernels to post-processing tools : such as ParaView throw MPI communication channel and follow in live the computation with no disk IO.

- Eye Dome Lighting

A non-photorealistic shading technique designed at EDF (France) to improve depth perception in scientific visualization images.

It relies on efficient post-processing passes implemented on the GPU with GLSL shaders in order to achieve interactive rendering.

- CoProcessingPlugin

Adds extensions to enable exporting state files that can be used by ParaView CoProcessing library.

- Force Time

Overrides the VTK time requests. This can be used to create complex animation with different datasets following independent times evolutions.

Note: As this filter overrides the time requests, time-aware filters such as PathLines or PlotOverTime will not behave correctly if they are inserted in a pipeline after this filter.

- H5PartReader

The H5Part Reader plugin adds support for reading particle datasets stored in H5Part format. H5Part is a simple wrapper around the HDF5 library and provides a number of convenience functions to manage time steps and access field arrays. The reader supports parallel reading of data using hyperslabs, when used with ParaView compiled with MPI and HDF5 with parallel IO enabled, the reader automatically uses hyperslabs to read portions of data on each process.

- Nifti

Reads time varying volumetric/image data from ANALYZE. Supports single (.nii) and dual (.img & .hdr) file storage including zlib compression.

Able to write out ascii or binary files.

- Manta View

A view that uses the University of Utah's Manta Real Time Ray Tracer instead of OpenGL for rendering surfaces

Kitware Source Article <sup>[1]</sup>

- Moments

Contains a set of filters that helps analysis of flux and circulation fields. Flux fields are defined on 2D cells and describe flows through the area of the cell. Circulation fields are defined on 1D cells and describes flows in the direction of the cell.

- PointSprite

Adds a renderer for point geometry - in particular particle based datasets (though any point based data may be rendered using the provided painter). The plugin permits 3 modes of rendering, which are (in increasing order

of complexity), Simple points, texture mapped sprites, and GPU raytraced spheres. The simple point mode allows the user to select a scalar array for the opacity on a per point basis. The texture mode adds support for opacity and radius per point (particle) which is drawn using a user supplied texture (a sphere is provided by default). The GPU mode differs by evaluating a quadric ray/sphere intersection that allow objects to intersect correctly rather than ‘popping’ in and out of view as sprites do. Transfer function editors can be used to map radius/opacity values if simple non-linear lookups are required

- **SierraPlotTools**

Adds toolbar buttons to carry out convenience macros such as toggling the background color between white and black, switching between surface and surface with edges display mode, and opening a wizard for generating plots of global, nodal, or element variables over time. This plugin works on Exodus data only

- **SLACTools**

An extension that streamline ParaView's user interface for Stanford Linear Accelerator users.

- **Streaming View**

Views that render in many streamed passes to reduce memory footprint and provide multiresolution rendering. This plugin replaces the Streaming and Adaptive ParaView derived applications.

Kitware Source Article <sup>[2]</sup>

- **SurfaceLIC**

Adds support for Line Integral Convolution over arbitrary surfaces.

- **Prism Plugin**

A Sandia contributed plugin for verification and debugging of simulation results. The plugin reads SESAME files to determine the inverse mapping from cartesian to phase space and allows users to visually debug simulations. Sesame files hold material model data (Equations of State, Opacities, Conductivities) in a tabular format as described by [http://t1web.lanl.gov/doc/SESAME\\_3Ddatabase\\_1992.html](http://t1web.lanl.gov/doc/SESAME_3Ddatabase_1992.html).

- **pvblot**

Implementation of a Sandia National Labs scripting language on top of ParaView.

- **VisTrails**

The VisTrails plugin for ParaView incorporates the provenance management capabilities of VisTrails into ParaView.

All of the actions a user performs while building and modifying a pipeline in ParaView are captured by the plugin. This allows navigation of all of the pipeline versions that have previously been explored.

For more information about this plugin see [3]

## References

- [1] <http://www.kitware.com/products/html/RenderedRealismAtNearlyRealTimeRates.html>
- [2] <http://www.kitware.com/products/html/MultiResolutionStreamingInVTKAndParaView.html>
- [3] [http://www.vistrails.org/index.php/ParaView\\_Plugin](http://www.vistrails.org/index.php/ParaView_Plugin)

# Loading Plugins

## Loading Plugins

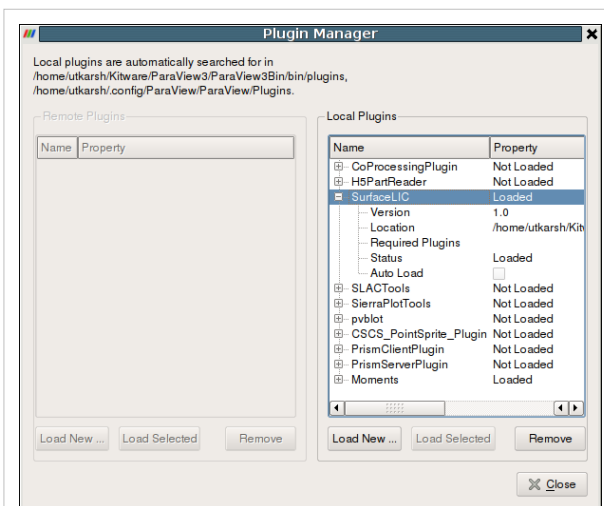
There are three ways for loading plugins:

- Using the GUI (**Plugin Manager**)

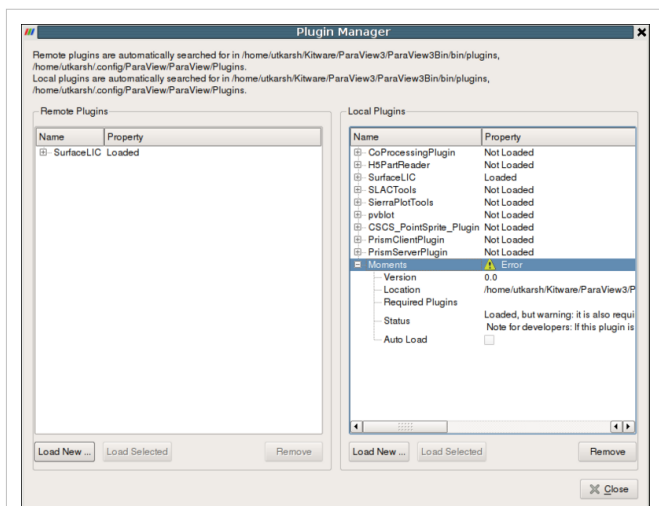
Plugins can be loaded into ParaView using the **Plugin Manager** accessible from **Tools | Manage Plugins/Extensions** menu. The Plugin Manager has two sections for loading local plugins and remote plugins (enabled only when connected to a server). To load a plugin on the local as well as remote side, simply browse to the plugin shared library. If the loading is successful, the plugin will appear in the list of loaded plugins. The Plugin manager also lists the paths it searched to load plugins automatically.

The Plugin Manager remembers all loaded plugins, so next time to load the plugin, simply locate it in the list and click "Load Selected" button.

You can set up ParaView to automatically load the plugin at startup (in case of client-side plugins) or on connecting to the server (in case of server-side plugins) by checking the "Auto Load" checkbox on a loaded plugin.



**Figure 1:** Plugin Manager when not connected to a remote server, showing loaded plugins on the local site.



**Figure 2:** Plugin Manager when connected to a server showing loaded plugins on the local as well as remote sites.

- Using environment variable (Auto-loading plugins)

If one wants ParaView to automatically load a set of plugins on startup, one can use the **PV\_PLUGIN\_PATH** environment variable. **PV\_PLUGIN\_PATH** can be used to list a set of directories (separated by colon (:) or semi-colon (;)) which ParaView will search on startup to load plugins. This environment variable needs to be set on both the client node to load local plugins as well as the remote server to load remote plugins. Note that plugins in **PV\_PLUGIN\_PATH** are always auto-loaded irrespective of the status of the "Auto Load" checkbox in the Plugin Manager.

- Placing the plugins in a recognized location. Recognized locations are:
  - A plugins subdirectory beneath the directory containing the paraview client or server executables. This can be a system-wide location if installed as such.
  - A Plugins subdirectory in the user's home area. On Unix/Linux/Mac, \$HOME/.config/ParaView/ParaView<version>/Plugins. On Windows

%APPDATA%\ParaView\ParaView<version>\Plugins.

---

# Appendix

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## Command Line Arguments

---

### Command-Line Arguments and Environment Variables

The following is a list of options available when running ParaView from the command line. When two options are listed, separated by a comma, either of them can be used to achieve the specified result. Unless otherwise specified, all command-line options are used on the client program. Following the list of command-line options is a set of environment variables ParaView recognizes.

#### General Options

- `--data` : Load the specified data file into ParaView (`--data=data_file`).
- `--disable-registry`, `-dr` : Launch ParaView in its default state; do not load any preferences saved in ParaView's registry file.
- `--help`, `/?` : Display a list of the command-line arguments and their descriptions.
- `--version`, `-V` : Display ParaView's version number, and then exit.

#### Client-Server Options

- `--server`, `-s` : Tell the client process where to connect to the server. The default is `--server=localhost`. This command-line option is used on the client.
  - `--client-host`, `-ch` : Tell the server process(es) where to connect to the client. The default is `--client-host=localhost`. This command-line option is used on the server(s).
  - `--server-port`, `-sp` : Specify the port to use in establishing a connection between the client and the server. The default is `--server-port=11111`. If used, this argument must be specified on both the client and the server command lines, and the port numbers must match.
  - `--data-server-port`, `-dsp` : Specify the port to use in establishing a connection between the client and the data server. The default is `--data-server-port=11111`. If used, this argument must be specified on both the client and the data server command lines, and the port numbers must match.
  - `--render-server-port`, `-rsp` : Specify the port to use in establishing a connection between the client and the render server. The default is `--render-server-port=22221`. If used, this argument must be specified on both the client and the render server command lines, and the port numbers must match.
  - `--reverse-connection`, `-rc` : Cause the data and render servers to connect to the client. When using this option, the client, data server, and render server (if used) must be started with this command-line argument, and you should start the client before starting either server.
  - `--connect-id` : Using a connect ID is a security feature in client-server mode. To use this feature, pass this command-line argument with the same ID number to the client and server(s). If you do not pass the same ID number, the server(s) will be shut down.
  - `--machines`, `-m` : Use this command-line argument to pass in the network configuration file for the render server. See section Error: Reference source not found for a description of this file.
-

## Rendering Options

- `--stereo` : Enable stereo rendering in ParaView.
- `--stereo-type`: Set the stereo rendering type. Options are:
  - Crystal Eyes
  - Red-Blue
  - Interlaced ( Default )
  - Dresden
  - Anaglyph
  - Checkerboard
- `--tile-dimensions-x, -tdx` : Specify the number of tiles in the horizontal direction of the tiled display (`--tile-dimensions-x=number_of_tiles`). This value defaults to 0. To use this option, you must be running in client/server or client/data server/render server mode, and this option must be specified on the client command line. Setting this option to a value greater than 0 will enable the tiled display. If this argument is set to a value greater than 0 and `-tdy` (see below) is not set, then `-tdy` will default to 1.
- `--tile-dimensions-y, -tdy` : Specify the number of tiles in the vertical direction of the tiled display (`--tile-dimensions-y=number_of_tiles`). This value defaults to 0. To use this option, you must be running in client/server or client/data server/render server mode, and this option must be specified on the client command line. Setting this option to a value greater than 0 will enable the tiled display. If this argument is set to a value greater than 0 and `-tdx` (see above) is not set, then `-tdx` will default to 1.
- `--tile-mullion-x, -tmx` : Specify the spacing (in pixels) between columns in tiled display images.
- `--tile-mullion-y, -tmy` : Specify the spacing (in pixels) between rows in tiled display images.
- `--use-offscreen-rendering` : Use offscreen rendering on the satellite processes. On unix platforms, software rendering or mangled Mesa must be used with this option.
- `--disable-composite, -dc` : Use this command-line option if the data server does not have rendering resources and you are not using a render server. All the rendering will then be done on the client.

## Environment Variables

In addition to the command-line options previously listed, ParaView also recognizes the following environment variables.

- `PV_DISABLE_COMPOSITE_INTERRUPTS` If this variable is set to 1, it is not possible to interrupt the compositing of images in parallel rendering. Otherwise it is interruptible through mouse interaction.
- `PV_ICET_WINDOW_BORDERS` Setting this variable to 1 when running ParaView in tiled display mode using IceT causes the render window for each tile to be the same size as the display area in ParaView's main application window. (Normally each render window fills the whole screen when tiled display mode is used.) This feature is sometimes useful when debugging ParaView.
- `PV_PLUGIN_PATH` If you have shared libraries containing plugins you wish to load in ParaView at startup, set this environment variable to the path for these libraries.
- `PV_SOFTWARE_RENDERING` This environment variable has the same effect as setting both the `--use-software-rendering` and `--use-satellite-software` environment variables.
- `VTK_CLIENT_SERVER_LOG` If set to 1, a log file will be created for the ParaView client, server, and render server. The log files will contain any client-server streams sent to the corresponding client, server, or render server.

## Executable help

paraview --help

```
--connect-id=opt      Set the ID of the server and client to make sure
they match.

--cslog=opt           ClientServerStream log file.

--data=opt            Load the specified data. To specify file series
replace the numeral with a '.' eg. my0.vtk, my1.vtk...myN.vtk becomes
my..vtk

--data-directory=opt  Set the data directory where test-case data are.

--disable-light-kit    When present, disables light kit by default.
Useful for dashboard tests.

--disable-registry
-dr                  Do not use registry when running ParaView (for
testing).

--exit               Exit application when testing is done. Use for
testing.

--help
/?                  Displays available command line arguments.

--machines=opt
-m=opt              Specify the network configurations file for the
render server.

--script=opt         Set a python script to be evaluated on startup.

--server=opt
-s=opt              Set the name of the server resource to connect
with when the client starts.

--state=opt          Load the specified statefile (.pvsm).

--stereo             Tell the application to enable stereo rendering
(only when running on a single process).

--stereo-type=opt     Specify the stereo type. This valid only when
--stereo is specified. Possible values are "Crystal Eyes", "Red-Blue",
"Interlaced", "Dresden", "Anaglyph", "Checkerboard"

--test-baseline=opt   Add test baseline. Can be used multiple times to
```



```

specify multiple baselines for multiple tests, in order.

--test-directory=opt  Set the temporary directory where test-case
output will be stored.

--test-script=opt      Add test script. Can be used multiple times to
specify multiple tests.

--test-threshold=opt  Add test image threshold. Can be used multiple
times to specify multiple image thresholds for multiple tests in order.

--version
-V                      Give the version number and exit.

--vrpn                  Tell the application to use VRPN for head
tracking

--vrpn-address=opt      Specify the VRPN tracker name. This valid only
when --vrpn is specified. Examples: "Tracker0@localhost",
"Head0@localhost"Please check VRPN configuration file

--vrui                  Tell the application to use VRUI for head
tracking

--vrui-address=opt      Specify the VRUI host name.

```

### **pvbatch --help**

```

--cslog=opt             ClientServerStream log file.

--help
/?                      Displays available command line arguments.

--machines=opt
-m=opt                  Specify the network configurations file for the render server.

--symmetric
-sym                    When specified, the python script is processed symmetrically on all processes.

--use-offscreen-rendering  Render offscreen on the satellite processes. This option only works with software rendering or mangled mesa on Unix.

--version
-V                      Give the version number and exit.

```

### **pvdatsarserver --help**

```

--client-host=opt
-ch=opt                  Tell the data|render server the host name of
the client, use with -rc.

```

```

--connect-id=opt      Set the ID of the server and client to make
sure they match.

--cslog=opt           ClientServerStream log file.

--data-server-port=opt
-dsp=opt              What port data server use to connect to the
client. (default 11111).

--help
/?                   Displays available command line arguments.

--machines=opt
-m=opt               Specify the network configurations file for
the render server.

--reverse-connection
-rc                  Have the server connect to the client.

--timeout=opt         Time (in minutes) since connecting with a
client after which the server may timeout. The client typically shows
warning messages before the server times out.

--version
-V                   Give the version number and exit.

```

### pvpython --help

```

--connect-id=opt      Set the ID of the server and client to make sure
they match.

--cslog=opt           ClientServerStream log file.

--data=opt            Load the specified data. To specify file series
replace the numeral with a '.' eg. my0.vtk, my1.vtk...myN.vtk becomes
my..vtk

--help
/?                   Displays available command line arguments.

--machines=opt
-m=opt               Specify the network configurations file for the
render server.

--state=opt           Load the specified statefile (.pvsm).

--stereo              Tell the application to enable stereo rendering

```

(only when running on a single process).

`--stereo-type=opt` Specify the stereo type. This valid only when `--stereo` is specified. Possible values are "Crystal Eyes", "Red-Blue", "Interlaced", "Dresden", "Anaglyph", "Checkerboard"

`--version`

`-V` Give the version number and exit.

### **pvrenderserver --help**

`--client-host=opt`

`-ch=opt` Tell the data|render server the host name of the client, use with `-rc`.

`--connect-id=opt` Set the ID of the server and client to make sure they match.

`--cslog=opt` ClientServerStream log file.

`--help`

`/?` Displays available command line arguments.

`--machines=opt`

`-m=opt` Specify the network configurations file for the render server.

`--render-server-port=opt`

`-rsp=opt` What port should the render server use to connect to the client. (default 22221).

`--reverse-connection`

`-rc` Have the server connect to the client.

`--tile-dimensions-x=opt`

`-tdx=opt` Size of tile display in the number of displays in each row of the display.

`--tile-dimensions-y=opt`

`-tdy=opt` Size of tile display in the number of displays in each column of the display.

`--tile-mullion-x=opt`

`-tmx=opt` Size of the gap between columns in the tile display, in Pixels.

`--tile-mullion-y=opt`

`-tmy=opt` Size of the gap between rows in the tile display, in Pixels.

`--use-offscreen-rendering` Render offscreen on the satellite processes. This option only works with software rendering or mangled mesa on Unix.

`--version`

`-V` Give the version number and exit.

### **pvserver --help**

```
--client-host=opt
  -ch=opt          Tell the data|render server the host name
of the client, use with -rc.

--connect-id=opt   Set the ID of the server and client to make
sure they match.

--cslog=opt        ClientServerStream log file.

--disable-composite
  -dc             Use this option when rendering resources
are not available on the server.

--help
  /?             Displays available command line arguments.

--machines=opt
  -m=opt          Specify the network configurations file for
the render server.

--reverse-connection
  -rc             Have the server connect to the client.

--server-port=opt
  -sp=opt         What port should the combined server use to
connect to the client. (default 11111).

--tile-dimensions-x=opt
  -tdx=opt        Size of tile display in the number of
displays in each row of the display.

--tile-dimensions-y=opt
  -tdy=opt        Size of tile display in the number of
displays in each column of the display.

--tile-mullion-x=opt
  -tmx=opt        Size of the gap between columns in the tile
display, in Pixels.

--tile-mullion-y=opt
  -tmy=opt        Size of the gap between rows in the tile
display, in Pixels.

--timeout=opt      Time (in minutes) since connecting with a
client after which the server may timeout. The client typically shows
warning messages before the server times out.
```

```
--use-offscreen-rendering  Render offscreen on the satellite  
processes. This option only works with software rendering or mangled  
mesa on Unix.
```

```
--version
```

-V Give the version number and exit.

## Application Settings

---

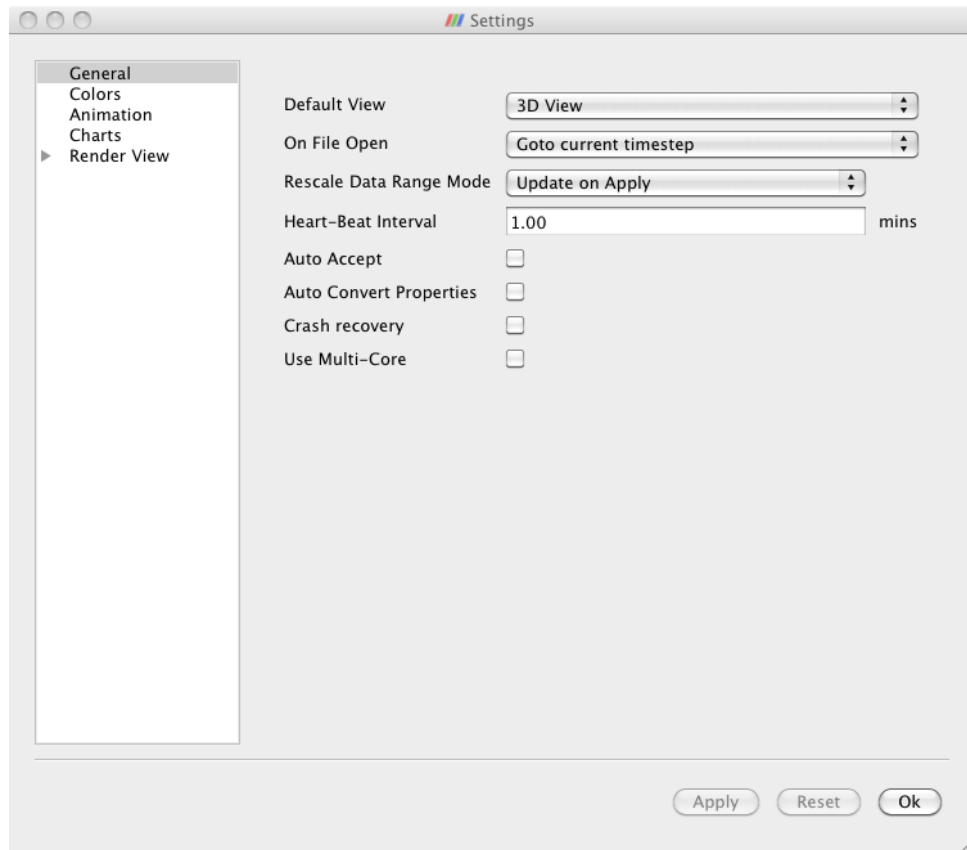
### Application Settings

Application wide settings are saved in between sessions so that every time you start ParaView on the same machine you get the same behaviors. The settings are stored in an XML file that resides at **%APPDATA%\ParaView\ParaViewVERSION.ini** on Windows and **~/.config/ParaView/ParaViewVERSION.ini** on all other platforms. If you need to restore ParaView's default settings you can either move this file or run with the **--disable-registry** argument.

These settings are managed via the settings dialog. On Macintosh this is reached from the **Preferences** entry in the **ParaView** menu. On Unix, Linux and Windows systems it is reached from the **Settings** entry on the **Edit** menu. These settings are program wide and are not to be confused with the View settings found under the **Edit** menu. Those settings are specific to each View window and are discussed in the Displaying Data <sup>[1]</sup> chapter.

The following figures show the various pages within the settings dialog. Choose from among them by clicking on the page list toward the left. Click on the arrow to expose the three sub pages in the **Render View** settings category. Some plugins add their own pages to the dialog but we do not discuss them here and instead refer you to any documentation that comes along with the plugin. Do note though that most settings are documented within the application via tooltips.

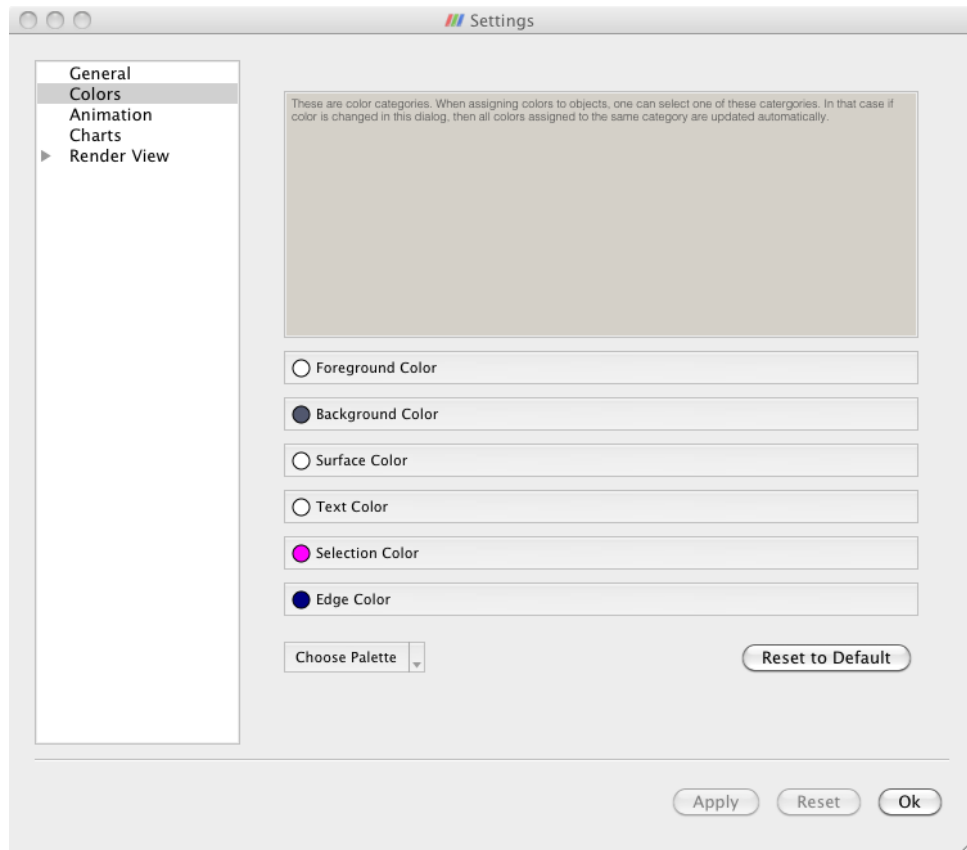
## General



**Figure 1. General Preferences**

- **Default View** : choose the View window type that ParaView initially shows when it starts up and connects to any server.
- **On File Open** : Control the timestep shown by default when ParaView opens a file with multiple time steps.
- **Rescale Data Range Mode** : Controls how color lookup tables are adjusted when data ranges change
- **Heart-Beat Interval** : Sets the length of time between keepalive messages that the client sends to the remote server which prevents the TCP connection between the two machines from shutting down.
- **Auto Accept** : Makes every change take effect immediately rather than requiring you to click the **Apply** button.
- **Auto Convert Properties** : Tells ParaView to do some data conversions automatically so that more filters are immediately applicable to the data you have at hand.
- **Crash Recovery** : Tells ParaView to save a record of each action you take so that you can easily recover your work if ParaView crashes mid-session.
- **Use Multi-Core** : Tells ParaView to spawn an MPI parallel server on the same machine that is running the GUI so as to better take advantage of multi-core machines on processing bound problems.

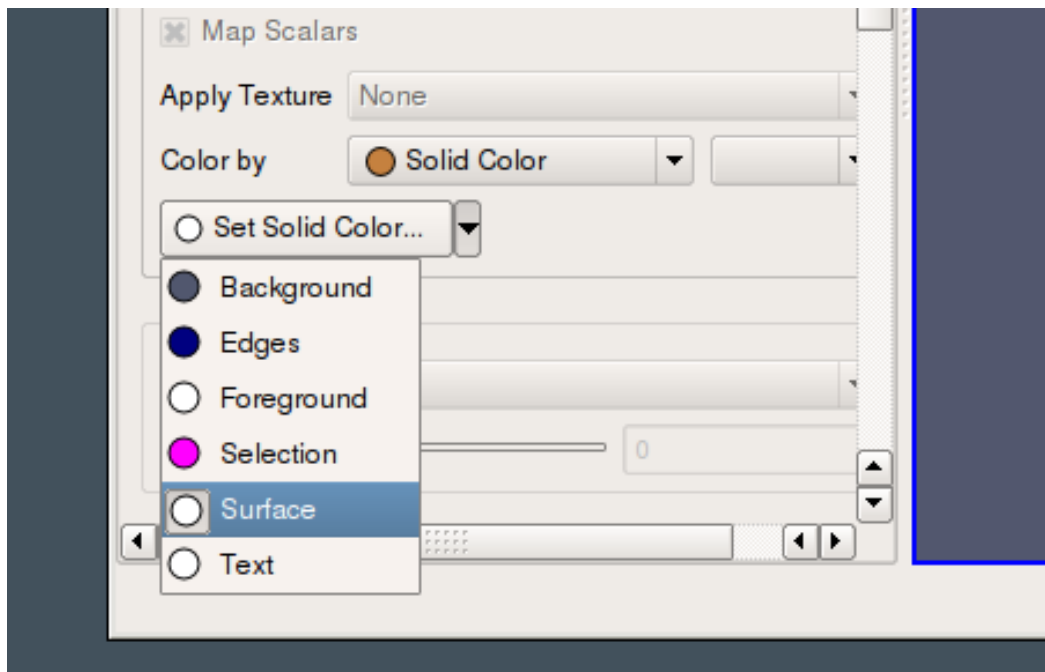
## Colors



**Figure 2. Color Preferences**

When setting up visualizations, most datasets in the visualization are typically colored using scalar colors. However, there may be items in the setup that are simply colored using a solid color including text and other annotations. Also by default ParaView uses a grey background, so the default colors chosen for objects tends to be setup so that it looks appropriate on the gray background. However, when saving a screenshot for printing, one would prefer to white background, for example. In that case, it can be tedious to go to every object and change its color to work with the new background color.

To make it easier to change colors on the fly, ParaView defines color categories. Users can choose to assign a color category to any object when specifying a color using the category menu on most color-chooser buttons, as shown below.



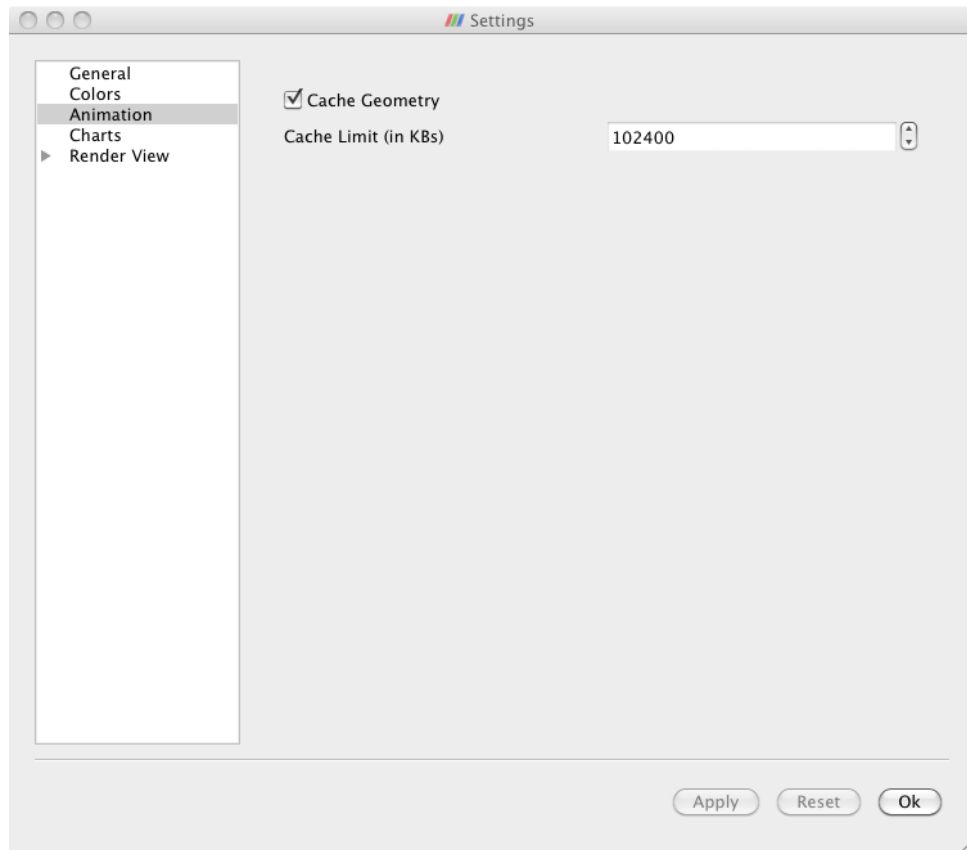
*Color Category Menu on the Display Tab*

Following associations are setup by default:

- Foreground color: outlines and wireframes
- Background color: background color in render views
- Surface color: solid color to surfaces
- Text color: color to 2D text
- Selection color: color used to show highlighted/selected elements in render view.
- Edge color: used for edges in **Surface With Edges** mode.



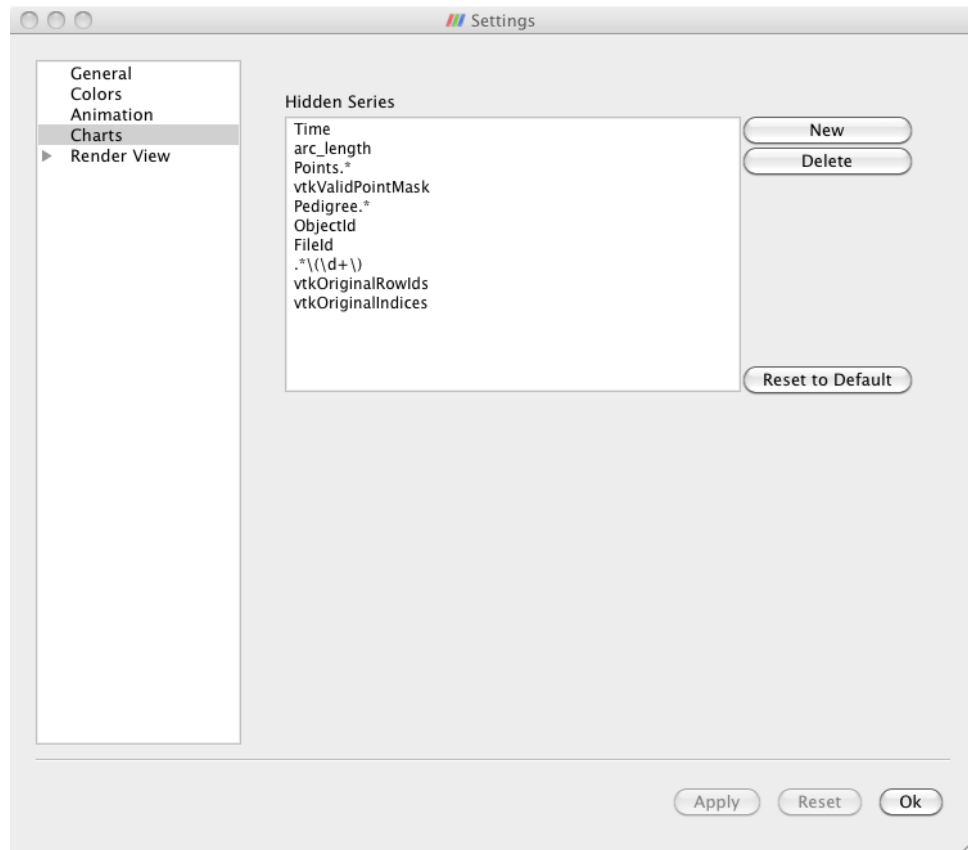
## Animation



**Figure 3. Animation Preferences**

- **Cache Geometry** : When checked this allows ParaView to save and later reuse rather than regenerate visible geometry during animation play back to speed up replay.
- **Cache Limit** : This is the maximum size of the animation playback geometry cache (on each node when in parallel).

## Charts



**Figure 4. Chart Preferences**

- **Hidden Series** : Allows you to specify a set of regular expressions that are compared against array names so that various bookkeeping arrays are not drawn in 2D charts and thus distract from the meaningful ones.

## Render View General

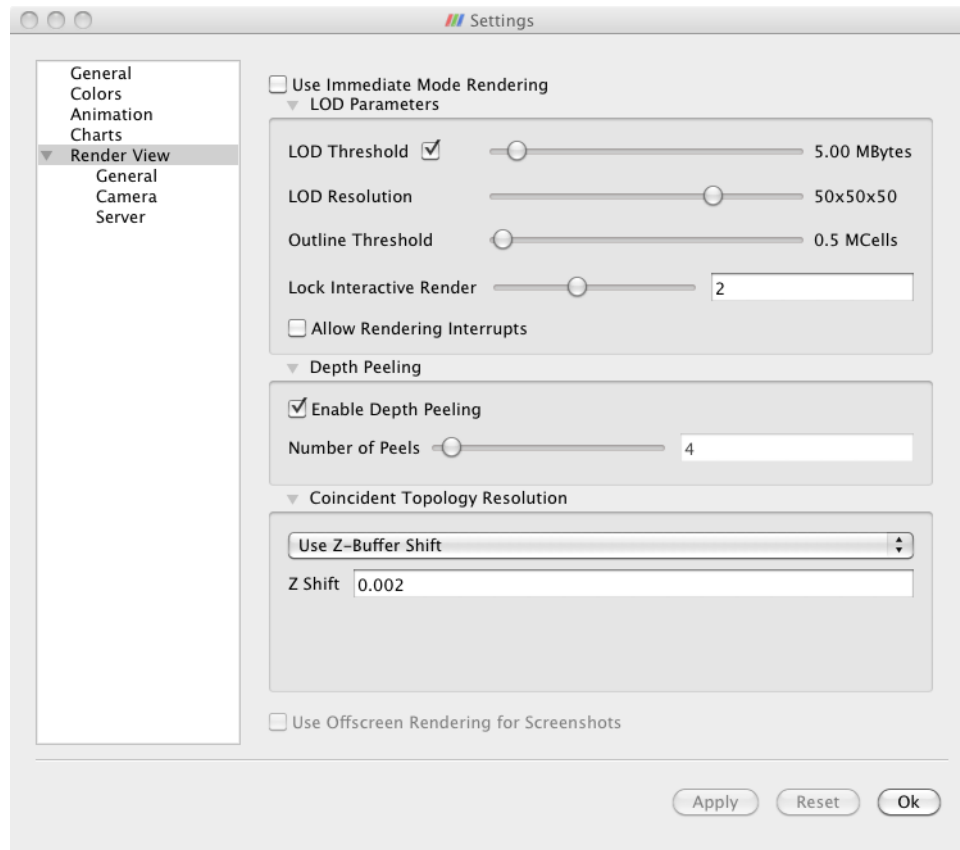


Figure 5. General Render View Preferences

- **Use Immediate Mode Rendering** : Controls whether OpenGL display lists are used or not which are generally faster only when rendering relatively small amounts of geometry.

The **LOD Parameters** settings control if, when and to what extent ParaView will down sample the data it draws while you move the mouse so as to maintain interactivity. The algorithm by which ParaView down samples the data is known as quadric clustering.

- **LOD Threshold** : Rendered data that is smaller than the specified size is not down sampled.
- **LOD Resolution** : When data is downsampled, this controls how coarsely.
- **Outline threshold** : Data that is larger than this threshold is drawn only as a bounding box while you move the camera.
- **Lock Interactive Render** : Controls how long ParaView waits after you have released the mouse and finished moving the camera before it returns to a full resolution render.
- **Allow Rendering Interrupts** : makes the drawing process interruptable so that you can move the camera immediately, without having to wait for a potentially slow full resolution render to complete.

The **Depth Peeling** settings control the algorithm that ParaView uses (given a sufficiently capable GPU) to draw translucent geometry correctly.

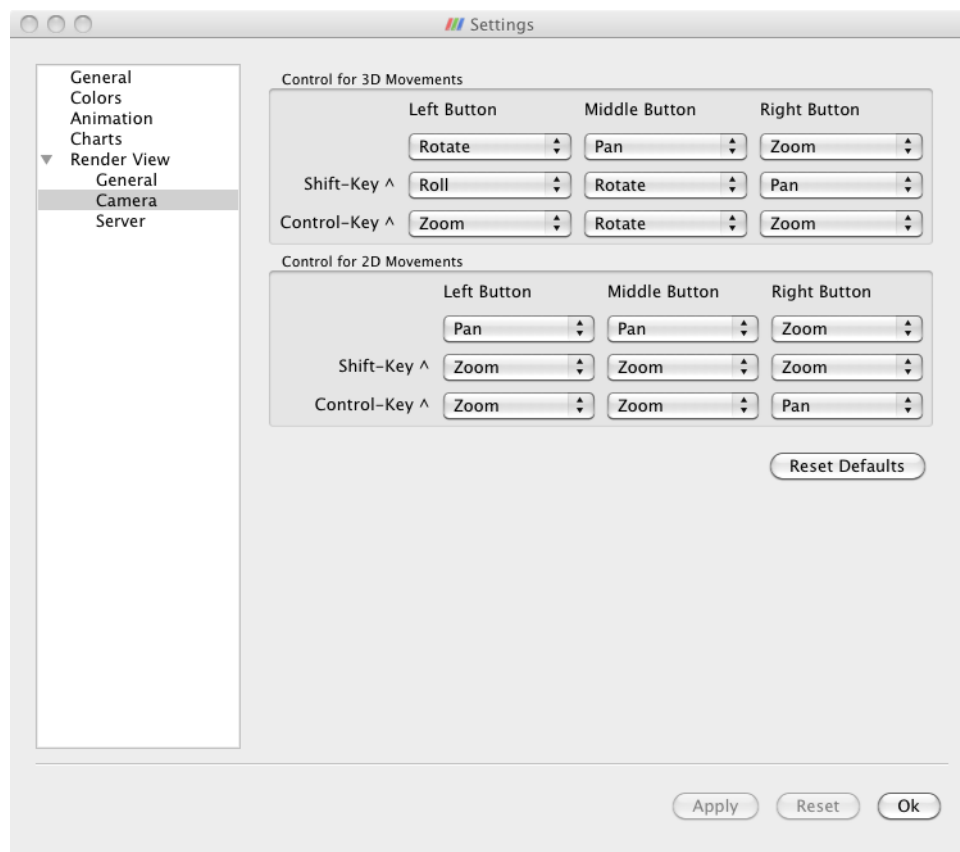
- **Enable Depth Peeling** : Controls whether ParaView will try to use the (potentially slow) depth peeling algorithm at all.
- **Number of Peels** : Controls the number of passes in the rendering algorithm. Higher produces quality images but increases rendering time.

The **Coincident Topology Resolution** settings control the method that ParaView uses to draw co-planar polygons in a non conflicting way (to avoid z-tearing in graphics terminology). Changes to these parameters do not take effect until ParaView is restarted.

The pull down menu allows you to choose between various algorithms that deal with z-tearing.

- **Do Nothing** : does not attempt to resolve overlapping geometry
- **Use Z-Buffer Shift** : adjusts OpenGL's Z direction scaling to draw one object slightly in front of the other. The **Z Shift** property controls the amount of offset.
- **Use Polygon Offset** : adjusts polygon locations (by calling `glPolygonOffset()`) for the same purpose. For an explanation of the **Factor** and **Units** parameters, see <http://www.opengl.org/resources/faq/technical/polygonoffset.htm>. The **Offset Faces** parameters controls whether filled polygons are moved rather than the points and edges.
- **Use Offscreen Rendering for Screenshots** : This checkbox is enabled only when offscreen rendering is available. It tells ParaView to generate images that are being captured as screenshots in offscreen contexts. On some platforms this is necessary to avoid accidentally capturing overlapping pixels (such as a mouse pointer) from the operating system.

## Render View Camera



**Figure 6. Camera Preferences**

These settings allow you to assign particular key/mouse press combinations to various camera moving controls. There are two sets of settings. **Control for 3D Movements** pertain to the 3D View. **Control for 2D Movements** pertain to the 2D View.

In the 3D View controls you can assign the following motions to any combination of left, middle or right mouse button press and no, shift or control key press.

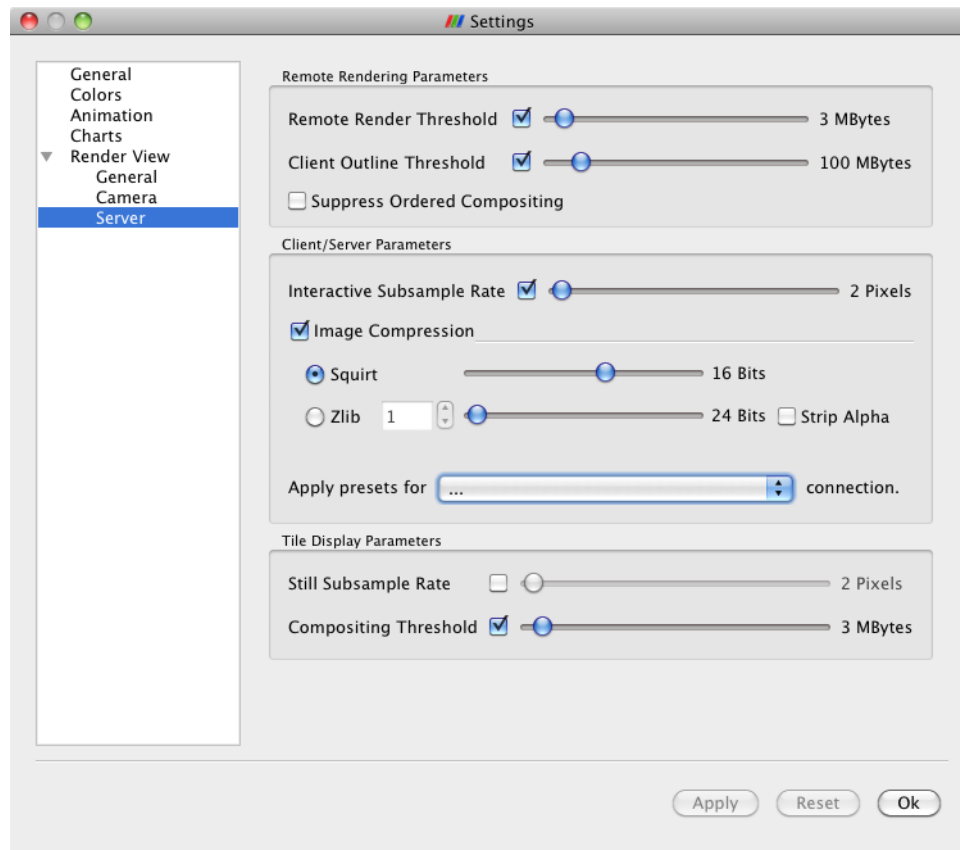
- **Pan**
- **Roll**
- **Rotate**
- **Multi-Rotate**

- **Zoom**

Rotation is not possible in the 2D View, which lacks a down direction to move about. So in this area you can only choose from **Pan** and **Zoom**.

Clicking **Reset Defaults** restores the button assignments to the factory settings.

## Render View Server



**Figure 7. Server Preferences**

These settings control how and where rendering happens when you are connected to a remote server.

The **Remote Rendering Parameters** are controls that generally affect the breakup of rendering work between the remote server and the local client machine.

- **Remote Render Theshold** : geometry that is smaller than the set value will be delivered to the client for local rendering. Data that is larger than the set value will be rendered remotely and pixels will be sent instead.
- **Client Outline Threshold** : while the camera is being manipulated and when geometry is being rendered locally, or in tiled display mode when it is being rendered both locally and remotely, data that is larger than this threshold will be rendered as a bounding box to maintain interactivity
- **Suppress Ordered Compositing** : This setting turns of the relatively slow transfer of data that must happen in parallel in order to render translucent geometry (as in volume rendering for instance) correctly.

The **Client/Server Parameters** section has controls that affect how pixels are transfered between the render server and the client while the data is being rendered remotely and images rather than geometry are being delivered to the client. The **Apply presets for ...** menu allows you to choose from a set of default settings which are generally well suited to particular network connection bandwidths.

- **Interactive Subsample Rate** : This setting controls how coarsely the image is downsampled to maintain interactivity while the camera is being manipulated.

- **Image Compression** : Enables compression of images during interaction. You can choose from either of the two image compression algorithms:
  - **Squirt** : Chooses IceT's native squirt image compression algorithm. The slider to the right controls the colorspace fidelity of the compression and conversely the interactive responsiveness.
  - **Zlib** : Chooses the ZLib library's compression algorithm instead. The spin box to the right specifies a compression level (1 through 9). The slider controls the colorspace fidelity. The **Strip Alpha** checkbox removed the Opacity channel (if any) from the shipped images.

The **Tile Display Parameters** section contains controls that take effect when doing tiled display wall rendering.

- **Still Subsample Rate** : This puts a screen space coarseness limit on the normally full resolution rendering that happens once you release the mouse after moving the camera. Image compositing time can prevent interactivity when working with large displays. This limit allows you to preview your results quickly and then choose to render truly full resolution results only once when you are satisfied with the preview.
- **Compositing Threshold** : This threshold allows you to get faster rendering times in the circumstance that you are visualizing relatively small data sets on large tiled displays. When the geometry size is below the threshold, the geometry is broadcast to all nodes in the render server. After this broadcast the relatively slow tiled image compositing algorithm is entirely omitted.

## List of Readers

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### AVS UCD Reader

Read binary or ASCII files stored in AVS UCD format.

The AVS UCD reader reads binary or ASCII files stored in AVS UCD format.

The default file extension is .inp. The output of this reader is unstructured grid. This supports reading a file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Names</b><br>( <i>FileNames</i> )          | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property contains a list of the cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property contains a list of the point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## BYU Reader

Read Movie.BYU files to produce polygonal data.

The BYU reader reads data stored in Movie.BYU format. The expected file extension is .g. The datasets resulting from reading these files are polygonal.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the BYU reader. |                  |              |

## COSMO Reader

Read a cosmology file into a vtkUnstructuredGrid.

The Cosmology reader reads a binary file of particle location, velocity, and id creating a vtkUnstructuredGrid. The default file extension is .cosmo. Reads LANL Cosmo format or Gadget format.

| Property  | Description  | Default Value(s) | Restrictions                                  |
|---|--|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                     | The list of files to be read by the reader.  |                  |   |
| <b>Cosmo Format</b><br>( <i>CosmoFormat</i> )                               | If checked, the data will be read as the LANL Cosmo format. If unchecked, the data will be read as Gadget-2 format.  | 1                | Only the values 0 and 1 are accepted.         |
| <b>overlap (shared point/ghost cell gap distance)</b><br>( <i>Overlap</i> ) | The space (in rL units) to extend processor particle ownership for ghost particles/cells. Needed for correct halo calculation when halos cross processor boundaries in parallel computation. | 5                | The value must be greater than or equal to 0. |
| <b>rL (physical box side length)</b><br>( <i>RL</i> )                       | The box side length used to wrap particles around if they exceed rL (or less than 0) in any dimension.   | 100              | The value must be greater than or equal to 0. |

## CSV Reader

Read a comma-separated values file into a 1D rectilinear grid.

The CSV reader reads a comma-separated values file into a 1D rectilinear grid. If the file was saved using the CSVWriter, then the rectilinear grid's points and point data can be restored as well as the cell data. Otherwise all the data in the CSV file is treated as cell data. The default file extension is .csv.

This can read file series as well.

| Property   | Description  | Default Value(s) | Restrictions                          |
|--|--|------------------|---------------------------------------|
| <b>File Name</b><br>( <i>FileName</i> )                                      | The list of files to be read by the reader. Each file is expected to be a csv file. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step.   |                  |                                       |
| <b>Detect Numeric Columns</b><br>( <i>DetectNumericColumns</i> )             | When set to true, the reader will detect numeric columns and create vtkDoubleArray or vtkIntArray for those instead of vtkStringArray.   | 1                | Only the values 0 and 1 are accepted. |
| <b>Field Delimiter Characters</b><br>( <i>FieldDelimiterCharacters</i> )     | This property lists the characters that may be used to separate fields. For example, a value of "," indicates a comma-separated value file. A value of ";;" indicates that columns may be separated by a period, colon or semicolon. The order of the characters in the text string does not matter. | ,                |                                       |
| <b>Have Headers</b><br>( <i>HaveHeaders</i> )                                | If the value of this property is 1, treat the first line of the file as headers.   | 1                | Only the values 0 and 1 are accepted. |
| <b>Merge Consecutive Delimiters</b><br>( <i>MergeConsecutiveDelimiters</i> ) | Whether to merge successive delimiters. Use this if (for example) your fields are separated by spaces but you don't know exactly how many.   | 0                | Only the values 0 and 1 are accepted. |
| <b>Use String Delimiter</b><br>( <i>UseStringDelimiter</i> )                 | This property indicates whether to use the string delimiter.   | 1                | Only the values 0 and 1 are accepted. |

## DEM Reader

Read a DEM (Digital Elevation Model) file.

The DEM reader reads Digital Elevation Model files containing elevation values derived from the U. S. Geologic Survey. The default file extension is .dem. This reader produces uniform rectilinear (image/volume) data output.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the DEM (Digital Elevation Map) reader. |                  |              |

## EnSight Master Server Reader

Read files in EnSight's Master Server format.

The EnSight Master Server reader reads EnSight's parallel files. The master file usually has a .sos extension and may point to multiple .case files. The output is a multi-block dataset.



| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Byte Order</b><br>( <i>ByteOrder</i> )          | This property indicates the byte order of the binary file(s).                           | 0                | The value must be one of the following: BigEndian (0), LittleEndian (1). |
| <b>Case File Name</b><br>( <i>CaseFileName</i> )   | This property specifies the name of the .sos file for the EnSight Master Server reader. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.                                 |                  | The list of array names is provided by the reader.                       |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.                                |                  | The list of array names is provided by the reader.                       |
| <b>Set Time Value</b><br>( <i>SetTimeValue</i> )   | This property indicates which time value to read.                                       | 0                |  |

## EnSight Reader

Read EnSight 6 and Gold files.

The EnSight reader reads files in the format produced by CEI's EnSight. EnSight 6 and Gold files (both ASCII and binary) are supported. The default extension is .case. The output of this reader is a multi-block dataset.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>Case File Name</b><br>( <i>CaseFileName</i> )   | This property specifies the case file name for the EnSight reader. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.            |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.           |                  | The list of array names is provided by the reader. |

## Enzo Reader

Read multi-block dataset from an Enzo file.

This Enzo reader loads data stored in Enzo format. The output of this reader is a multi-block dataset.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Block Output Type</b><br>( <i>BlockOutputType</i> ) | The property indicates the output type of each block, either vtkImageData (by default) or vtkRectilinearGrid. | 0                | The value must be one of the following: Image Data (0), Rectilinear Grid (1). |
| <b>File Name</b><br>( <i>FileName</i> )                | This property specifies the file name for the Enzo reader.  |                  |   |
| <b>Load Particles</b><br>( <i>LoadParticles</i> )      | Load particles when on.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Max Level</b><br>( <i>MaxLevel</i> )                | Do not blocks above this level.   | 100              |   |

## ExodusIIReader

Read an Exodus II file to produce an unstructured grid.

The Exodus reader loads Exodus II files and produces an unstructured grid output. The default file extensions are .g, .e, .ex2, .ex2v2, .exo, .gen, .exoII, .exii, .0, .00, .000, and .0000. The file format is described fully at: <http://endo.sandia.gov/SEACAS/Documentation/exodusII.pdf>.

Each Exodus file contains a single set of points with 2-D or 3-D coordinates plus one or more blocks, sets, and maps. Block group elements (or their bounding edges or faces) of the same type together. Sets select subsets (across all the blocks in a file) of elements, sides of elements (which may be of mixed dimensionality), bounding faces of volumetric elements, or bounding edges of volumetric or areal elements. Each block or set may have multiple result variables, each of which defines a value per element, per timestep.

The elements (cells), faces of elements (when enumerated in face blocks), edges of elements (when enumerated in edge blocks), and nodes (points) in a file may be assigned an arbitrary integer number by an element map, face map, edge map, or node map, respectively. Usually, only a single map of each type exists and is employed to assign a unique global ID to entities across multiple files which partition a large mesh for a distributed-memory calculation. However here may be multiply maps of each type and there are no constraints which force the integers to be unique.

The connectivity of elements is constant across all of the timesteps in any single Exodus II file. However, multiple files which specify a single time-evolution of a mesh may be used to represent meshes which exhibit changes in connectivity infrequently.

| Property  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                           | This property specifies the file name for the Exodus reader.  |                  |  |
| <b>Use Meta File</b><br>( <i>UseMetaFile</i> )                    | This hidden property must always be set to 1 for this proxy to work.  | 0                | Only the values 0 and 1 are accepted.              |
| <b>Animate Vibrations</b><br>( <i>AnimateVibrations</i> )         | If this flag is on and HasModeShapes is also on, then this reader will report a continuous time range [0,1] and animate the displacements in a periodic sinusoid. If this flag is off and HasModeShapes is on, this reader ignores time. This flag has no effect if HasModeShapes is off.   | 1                | Only the values 0 and 1 are accepted.              |
| <b>Apply Displacements</b><br>( <i>ApplyDisplacements</i> )       | Geometric locations can include displacements. When this option is on, the nodal positions are 'displaced' by the standard exodus displacement vector. If displacements are turned 'off', the user can explicitly add them by applying a warp filter.   | 1                | Only the values 0 and 1 are accepted.              |
| <b>Displacement Magnitude</b><br>( <i>DisplacementMagnitude</i> ) | When displacements are being applied, they are scaled by this amount. Set to 1 for no scaling.  | 1                |  |
| <b>Edge Blocks</b><br>( <i>EdgeBlocks</i> )                       | Use this property to select which blocks of edges should be loaded. All edges in the same block will have the same interpolant (linear, quadratic, etc.) and have the same variables defined over them.   |                  | The list of array names is provided by the reader. |
| <b>Edge Map Array Status</b><br>( <i>EdgeMapArrayStatus</i> )     | Specify which maps defined over edges in the mesh should be loaded. An edge map is a renumbering of all the edges (across all the edge blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files.   |                  | The list of array names is provided by the reader. |
| <b>Edge Set Array Status</b><br>( <i>EdgeSetArrayStatus</i> )     | An Exodus II file may define subsets of all the edges (across all edge blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per edge. The accompanying EdgeSetResultArrayStatus property specifies which variables defined over those sets should be loaded. |                  | The list of array names is provided by the reader. |

|  |   |   |  |
|--|---|---|--|
| <b>Edge Set Result Array Status</b><br>( <i>EdgeSetResultArrayStatus</i> )       | An Exodus II file may define subsets of all the edges (across all edge blocks) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per edge. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the EdgeSetArrayStatus property in order for the variable to appear.                          |   | The list of array names is provided by the reader. |
| <b>Edge Variables</b><br>( <i>EdgeVariables</i> )                                | Use this property to select which variables defined over edges should be loaded. Edges are oriented boundaries of one or more higher-dimensional cells also present in the mesh. Edge variables specify a single value per entry in some edge block.  |   | The list of array names is provided by the reader. |
| <b>Element Blocks</b><br>( <i>ElementBlocks</i> )                                | Use this property to select which blocks of elements (cells) should be loaded. All elements (cells) in the same block will have the same interpolant (tetrahedral, triangular, trilinear hexahedral, etc.) and have the same variables defined over them.   |   |  |
| <b>Element Map Array Status</b><br>( <i>ElementMapArrayStatus</i> )              | Specify which maps defined over elements (cells) in the mesh should be loaded. An element map is a renumbering of all the elements (across all the element blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files. These are sometimes called new-style maps because they are specified differently than in older Exodus files, which only allowed a single element map. |   | The list of array names is provided by the reader. |
| <b>Element Set Array Status</b><br>( <i>ElementSetArrayStatus</i> )              | An Exodus II file may define subsets of all the elements (across all element blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as feature classifiers, may then be defined over these sets by specifying a single number per element. The accompanying ElementSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                 |   | The list of array names is provided by the reader. |
| <b>Element Set Result Array Status</b><br>( <i>ElementSetResultArrayStatus</i> ) | An Exodus II file may define subsets of all the elements (across all element blocks) in a file as sets in their own right. Variables, such as feature classifiers, may then be defined over these sets by specifying a single number per element. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the ElementSetArrayStatus property in order for the variable to appear.              |   | The list of array names is provided by the reader. |
| <b>Element Variables</b><br>( <i>ElementVariables</i> )                          | Use this property to select which variables defined over elements (cells) should be loaded. Element variables specify a single value per entry in some element block.   |   | The list of array names is provided by the reader. |
| <b>Exodus Model Metadata</b><br>( <i>ExodusModelMetadata</i> )                   | Should metadata for the mesh be stored in the field data of the reader's output?  | 0 | Only the values 0 and 1 are accepted.              |
| <b>Face Blocks</b><br>( <i>FaceBlocks</i> )                                      | Use this property to select which blocks of faces should be loaded. All faces in the same block will have the same interpolant (triangular, bilinear quadrilateral, etc.) and have the same variables defined over them.  |   | The list of array names is provided by the reader. |
| <b>Face Map Array Status</b><br>( <i>FaceMapArrayStatus</i> )                    | Specify which maps defined over faces in the mesh should be loaded. A face map is a renumbering of all the faces (across all the face blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files.  |   | The list of array names is provided by the reader. |
| <b>Face Set Array Status</b><br>( <i>FaceSetArrayStatus</i> )                    | An Exodus II file may define subsets of all the faces (across all face blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per face. The accompanying FaceSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                             |   | The list of array names is provided by the reader. |

|   |  |       |  |
|---|--|-------|--|
| <b>Face Set Result Array Status</b><br>(FaceSetResultArrayStatus)         | An Exodus II file may define subsets of all the faces (across all face blocks) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per face. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the FaceSetArrayStatus property in order for the variable to appear. |       | The list of array names is provided by the reader. |
| <b>Face Variables</b><br>(FaceVariables)                                  | Use this property to select which variables defined over faces should be loaded. Faces are oriented boundaries of one or more higher-dimensional cells also present in the mesh. Face variables specify a single value per entry in some face block.   |       | The list of array names is provided by the reader. |
| <b>File Pattern</b><br>(FilePattern)                                      | This property contains a text string specifying the pattern used to load files. There are two parts to it: the prefix (%s) and the file number (%i). The format used is the same as that used for printf.  |       |  |
| <b>File Prefix</b><br>(FilePrefix)  | This property specifies the file root used with the file pattern to format a file name. (See the File Pattern property.)   |       |  |
| <b>File Range</b><br>(FileRange)  | This property controls the indices of the first and last files to be read.   | -1 -1 |  |
| <b>Generate Global Element Id Array</b><br>(GenerateGlobalElementIdArray) | Should the reader retrieve the first new-style element map, or if that is not present, the solitary old-style element map (which always exists but may be procedurally generated if it is not stored with the file)?   | 1     | Only the values 0 and 1 are accepted.              |
| <b>Generate Global Node Id Array</b><br>(GenerateGlobalNodeIdArray)       | Should the reader retrieve the first new-style node map, or if that is not present, the solitary old-style node map (which always exists but may be procedurally generated if it is not stored with the file)?   | 1     | Only the values 0 and 1 are accepted.              |
| <b>Generate Object Id Cell Array</b><br>(GenerateObjectIdCellArray)       | Specifies whether a variable named "ObjectId" should be created for each entry of each block and set being loaded and assigned the identifying number of the block or set, so that given a cell one can readily identify the block or set it came from. This is useful for coloring a mesh by block ID or set ID.  | 1     | Only the values 0 and 1 are accepted.              |
| <b>Global Variables</b><br>(GlobalVariables)                              | Use this property to select which global (i.e., per-mesh) variables should be loaded. Global variables are stored as arrays in the field data of the reader's output.  |       | The list of array names is provided by the reader. |
| <b>Has Mode Shapes</b><br>(HasModeShapes)                                 | Some simulations overload the Exodus time steps to represent mode shapes. In this case, it does not make sense to iterate over the "time steps", because they are not meant to be played in order. Rather, each represents the vibration at a different "mode." Setting the value of this property to 1 changes the semantics of the reader to not report the time steps to downstream filters.                                  | 0     | Only the values 0 and 1 are accepted.              |
| <b>Mode Shape</b><br>(ModeShape)  | Specify which mode shape to animate when HasModeShapes is on.  | 1     |  |
| <b>Node Map Array Status</b><br>(NodeMapArrayStatus)                      | Specify which maps defined over nodes (points) in the mesh should be loaded. A node map is a renumbering of all the points in a file from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files. These are sometimes called new-style maps because they are specified differently than in older Exodus files, which only allowed a single node map.                   |       | The list of array names is provided by the reader. |
| <b>Node Set Array Status</b><br>(NodeSetArrayStatus)                      | An Exodus II file may define subsets of all the nodes (points) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per node. This property specifies which of those sets should be loaded. The accompanying NodeSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                    |       | The list of array names is provided by the reader. |

|  |   |  |  |
|--|---|--|--|
| <b>Node Set Result Array Status</b><br>( <i>NodeSetResultArrayStatus</i> ) | An Exodus II file may define subsets of all the nodes (points) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per node. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the NodeSetArrayStatus property in order for the variable to appear.  |  | The list of array names is provided by the reader. |
| <b>Point Variables</b><br>( <i>PointVariables</i> )                        | Use this property to select which nodal (i.e., per-point) variables should be loaded.   |  | The list of array names is provided by the reader. |
| <b>Side Set Array Status</b><br>( <i>SideSetArrayStatus</i> )              | An Exodus II file may define subsets of all the of all the elements in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per side. For example, a hexahedron has 18 sides: 6 faces and 12 edges. Any of these sides may be individually called out in a set and assigned a result value. The accompanying SideSetResultArrayStatus property specifies which variables defined over those sets should be loaded.    |  | The list of array names is provided by the reader. |
| <b>Side Set Result Array Status</b><br>( <i>SideSetResultArrayStatus</i> ) | An Exodus II file may define subsets of all the of all the elements in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per side. For example, a hexahedron has 18 sides: 6 faces and 12 edges. Any of these sides may be individually called out in a set and assigned a result value. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the SideSetArrayStatus property in order for the variable to appear. |  | The list of array names is provided by the reader. |
| <b>XML File Name</b><br>( <i>XMLFileName</i> )                             | This property specifies the name of an XML file containing part and material descriptions and a mapping to element blocks.  |  |  |

## FLUENTReader

Reads a dataset in Fluent file format.

FLUENTReader creates an unstructured grid dataset. It reads .cas and .dat files stored in FLUENT native format.

| Property   | Description                                | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> ) | Select which cell-centered arrays to read. |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )          | Set the file name for the FLUENT reader.   |                  |  |

## Facet Reader

Read ASCII files stored in Facet format.

The Facet Reader reads files in Facet format: a simple ASCII file format listing point coordinates and connectivity between these points. The default file extension is .facet. The output of the Facet Reader is polygonal.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the Facet reader. |                  |              |

## Flash Reader

Read multi-block dataset from a Flash file.

This Flash reader loads data stored in Flash format. The output of this reader is a multi-block dataset.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Block Output Type</b><br>( <i>BlockOutputType</i> ) | The property indicates the output type of each block, either vtkImageData (by default) or vtkRectilinearGrid. | 0                | The value must be one of the following: Image Data (0), Rectilinear Grid (1). |
| <b>File Name</b><br>( <i>FileName</i> )                | This property specifies the file name for the Flash reader.   |                  |   |
| <b>Load Morton Curve</b><br>( <i>LoadMortonCurve</i> ) | Generate morton curve when on.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Load Particles</b><br>( <i>LoadParticles</i> )      | Load particles when on.   | 1                | Only the values 0 and 1 are accepted.   |

## Gaussian Cube Reader

Produce polygonal data by reading a Gaussian Cube file.

The Gaussian Cube reader produces polygonal data by reading data files produced by the Gaussian software package. The expected file extension is .cube.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>B Scale</b><br>( <i>BScale</i> )     | A scaling factor to compute bonds between non-hydrogen atoms        | 1                |              |
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the Gaussian Cube reader. |                  |              |
| <b>H BScale</b><br>( <i>HBScale</i> )   | A scaling factor to compute bonds with hydrogen atoms.              | 1                |              |

## Image Reader

Read raw regular rectilinear grid data from a file. The dimensions and type of the data must be specified.

The Image reader reads raw regular rectilinear grid (image/volume) data from a file. Because no metadata is provided, the user must specify information about the size, spacing, dimensionality, etc. about the dataset.

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Data Byte Order</b><br>( <i>DataByteOrder</i> ) | This property indicates the byte order of the binary file(s).  | 0                | The value must be one of the following: BigEndian (0), LittleEndian (1). |
| <b>Data Extent</b><br>( <i>DataExtent</i> )        | This property specifies the minimum and maximum index values of the data in each dimension (xmin, xmax, ymin, ymax, zmin, zmax). | 0 0 0 0 0        |  |
| <b>Data Origin</b><br>( <i>DataOrigin</i> )        | The coordinate contained in this property specifies the position of the point with index (0,0,0).                                | 0 0 0            |  |

|   |  |           |  |
|---|--|-----------|--|
| <b>Data Scalar Type</b><br>( <i>DataScalarType</i> )                      | The value of this property indicates the scalar type of the pixels/voxels in the file(s): short, int, float ...  | 4         | The value must be one of the following: char (2), unsigned char (3), short (4), unsigned short (5), int (6), unsigned int (7), long (8), unsigned long (9), float (10), double (11). |
| <b>Data Spacing</b><br>( <i>DataSpacing</i> )                             | This property specifies the size of a voxel in each dimension.   | 1 1 1     |  |
| <b>File Dimensionality</b><br>( <i>FileDimensionality</i> )               | This property indicates whether the file(s) in this dataset contain slices (2D) or volumes (3D).   | 3         | The value must be one of the following: 2 (2), 3 (3).  |
| <b>File Lower Left</b><br>( <i>FileLowerLeft</i> )                        | This property determines whether the data originates in the lower left corner (on) or the upper left corner (off). Most scientific data is written with a right-handed axes that originates in the lower left corner. However, several 2D image file formats write the image from the upper left corner.           | 1         | Only the values 0 and 1 are accepted.  |
| <b>File Pattern</b><br>( <i>FilePattern</i> )                             | The text string contained in the property specifies the format string to determine the file names necessary for reading this dataset. In creating the filenames, %s will be replaced by the prefix and %d by a digit which represents the slice number in Z. The format string is the same as that used by printf. | %s        |  |
| <b>File Prefix</b><br>( <i>FilePrefix</i> )                               | The text string contained in this property specifies the file prefix (directory plus common initial part of file name) for the raw binary uniform rectilinear grid dataset.  |           |  |
| <b>Number Of Scalar Components</b><br>( <i>NumberOfScalarComponents</i> ) | This property specifies the number of components the scalar value at each pixel or voxel has (e.g., RGB - 3 scalar components).  | 1         |  |
| <b>Scalar Array Name</b><br>( <i>ScalarArrayName</i> )                    | This property contains a text string listing a name to assign to the point-centered data array read.   | ImageFile |  |

## JPEG Series Reader

Read a series of JPEG files into an time sequence of image datas.

The JPEG series reader reads JPEG files. The output is a time sequence of uniform rectilinear (image/volume) dataset. The default file extension is .jpg or .jpeg.

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## LSDynReader

Read LS-Dyna databases (d3plot).

This reader reads LS-Dyna databases.

| Property  | Description  | Default Value(s) | Restrictions                                       |
|---|--|------------------|--|
| <b>Beam Arrays</b><br>( <i>BeamArrayStatus</i> )                | Select which beam arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Deformed Mesh</b><br>( <i>DeformedMesh</i> )                 | Should the mesh be deformed over time (if the Deflection node array is set to load)?             | 1                | Only the values 0 and 1 are accepted.              |
| <b>File Name</b><br>( <i>FileName</i> )                         | Set the file name for the LSDyna reader.   |                  |  |
| <b>Part Arrays</b><br>( <i>PartArrayStatus</i> )                | Select which part arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Particle Arrays</b><br>( <i>ParticleArrayStatus</i> )        | Select which particle arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> )              | Select which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Remove Deleted Cells</b><br>( <i>RemoveDeletedCells</i> )    | Should cells that have been deleted (failed structurally, for example) be removed from the mesh? | 1                | Only the values 0 and 1 are accepted.              |
| <b>Rigid Body Arrays</b><br>( <i>RigidBodyArrayStatus</i> )     | Select which rigid body arrays to load.  |                  | The list of array names is provided by the reader. |
| <b>Road Surface Arrays</b><br>( <i>RoadSurfaceArrayStatus</i> ) | Select which road surface arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Shell Arrays</b><br>( <i>ShellArrayStatus</i> )              | Select which shell arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Solid Arrays</b><br>( <i>SolidArrayStatus</i> )              | Select which solid arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Split by Material Id</b><br>( <i>SplitByMaterialId</i> )     | Should each material have its own mesh, or should there be one mesh for all materials?           | 0                | Only the values 0 and 1 are accepted.              |
| <b>Thick Shell Arrays</b><br>( <i>ThickShellArrayStatus</i> )   | Select which thick shell arrays to read.   |                  | The list of array names is provided by the reader. |

## Legacy VTK Reader

Read files stored in VTK's legacy file format.

The Legacy VTK reader loads files stored in VTK's legacy file format (before VTK 4.2, although still supported). The expected file extension is .vtk. The type of the dataset may be structured grid, uniform rectilinear grid (image/volume), non-uniform rectilinear grid, unstructured grid, or polygonal. This reader also supports file series.

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |



## MFIXReader

Reads a dataset in MFIX file format.

vtkMFIXReader creates an unstructured grid dataset. It reads a restart file and a set of sp files. The restart file contains the mesh information. MFIX meshes are either cylindrical or rectilinear, but this reader will convert them to an unstructured grid. The sp files contain transient data for the cells. Each sp file has one or more variables stored inside it.

| Property   | Description                                | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> ) | Select which cell-centered arrays to read. |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )          | Set the file name for the MFIX reader.     |                  |  |

## Meta File Series Reader

Read a series of meta images.

Read a series of meta images. The file extension is .mhd

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. Each file is expected to be in the meta format. The standard extension is .mhd. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## NetCDF MPAS reader

Read unstructured grid MPAS data from a file.

This reader reads unstructured grid MPAS data from a file. It creates a dual grid. It assumes the grid is in the global domain. By default, it creates a spherical view of vertical layer 0. It assumes it is ocean data. It gives several options to change the view to multilayer (all vertical layers will have a thickness determined by the value in the slider), lat/lon projection or atmospheric (vertical layers go out away from the center of the sphere, instead of down towards the center as they do for ocean data). It doesn't handle data in the rectangular domain. This is not a parallel reader. It expects one .nc file of data. It can display cell or vertex-centered data, but not edge data. When converted to the dual grid, cell-centered data becomes vertex-centered and vice-versa.

NOTES:

When using this reader, it is important that you remember to do the following:

1. When changing a selected variable, remember to select it also in the drop down box to color by. It doesn't color by that variable automatically
2. When selecting multilayer sphere view, start with layer thickness around 100,000
3. When selecting multilayer lat/lon view, start with layer thickness around 10
4. Always click the -Z orientation after making a switch from lat/lon to sphere, from single to multilayer or changing thickness.

5. Be conservative on the number of changes you make before hitting Apply, since there may be bugs in this reader. Just make one change and then hit Apply.

For problems, contact Christine Ahrens (cahrens@lanl.gov)

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>Cell Array Status</b><br>( <i>CellArrayStatus</i> )       | This property lists which NetCDF dual-grid cell variables to load.   |                  | The list of array names is provided by the reader. |
| <b>Center Longitude</b><br>( <i>CenterLon</i> )              | This property specifies where the center will be viewed for a lat/lon projection.  | 180              |  |
| <b>File Name</b><br>( <i>FileName</i> )                      | This property specifies the file name to read. It should be an MPAS format NetCDF file ending in .nc.  |                  |  |
| <b>Is Atmosphere</b><br>( <i>IsAtmosphere</i> )              | This property indicates whether data is atmospheric. Checking this ensures the vertical levels will go up away from the sphere instead of down towards the center.   | 0                | Only the values 0 and 1 are accepted.              |
| <b>Layer Thickness</b><br>( <i>LayerThickness</i> )          | This property specifies how thick the layer should be if viewing the data in multilayer view. Each layer corresponds to a vertical level. A good starting point is 100,000 for the spherical view and 10 for the lat/lon projection. Click on -Z after applying this change, since the scale may change drastically.   | 10               |  |
| <b>Point Array Status</b><br>( <i>PointArrayStatus</i> )     | This property lists which NetCDF dual-grid point variables to load.  |                  | The list of array names is provided by the reader. |
| <b>Project Lat Lon</b><br>( <i>ProjectLatLon</i> )           | This property indicates whether to view the data in the lat/lon projection.  | 0                | Only the values 0 and 1 are accepted.              |
| <b>Show Multilayer View</b><br>( <i>ShowMultilayerView</i> ) | This property indicates whether to show multiple layers in one view, with each vertical level having the same thickness, specified by the layer thickness slider. For ocean data, the layers correspond to data at vertical level whose number increases towards the center of the sphere. For atmospheric data, the layers correspond to data at vertical levels increasing away from the center. | 0                | Only the values 0 and 1 are accepted.              |
| <b>Vertical Level</b><br>( <i>VerticalLevel</i> )            | This property specifies which vertical level is viewed if not in multilayer view. Only the data for that vertical level will be viewed. The grid is essentially the same for each vertical level, however at some ocean levels, especially the lower ones, due to the topography, the grid becomes more sparse where there is land.  | 0                |  |

## NetCDF POP reader

Read raw regular rectilinear grid data from a file. The dimensions and type of the data must be specified.

The Image reader reads raw regular rectilinear grid (image/volume) data from a file. Because no metadata is provided, the user must specify information about the size, spacing, dimensionality, etc. about the dataset.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name to read.   |                  |   |
| <b>Stride</b><br>( <i>Stride</i> )                 | This property indicate the number of indices per dimension (i,j,k) to skip between each point included in this output. | 1 1 1            | The value must be greater than or equal to (1, 1, 1). |
| <b>Variables</b><br>( <i>VariableArrayStatus</i> ) | This property lists which netCDF variables to load.  |                  | The list of array names is provided by the reader.    |

## NetCDF Reader

Read regular arrays from netCDF files. Will also read any topological information specified by the COARDS and CF conventions.

Reads arrays from netCDF files into structured VTK data sets. In the absence of any other information, the files will be read as image data. This reader will also look for meta information specified by the CF convention that specify things like topology and time. This information can cause the output to be a nonuniform rectilinear grid or curvilinear (structured) grid. Details on the CF convention can be found at <http://cf-pcmdi.llnl.gov/>. It should be noted that the CF convention is a superset of the COARDS convention, so COARDS conventions will also be recognized.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                  | The name of the files to load.  |                  |   |
| <b>Dimensions</b><br>( <i>Dimensions</i> )                               | Load the grid with the given dimensions. Any arrays that conform to these dimensions will be loaded.  |                  | .   |
| <b>Output Type</b><br>( <i>OutputType</i> )                              | Specifies the type of data that the reader creates. If Automatic, the reader will use the most appropriate grid type for the data read. Note that not all grid types support all data. A warning is issued if a mismatch occurs.  | -1               | The value must be one of the following: Automatic (-1), Image (6), Rectilinear (3), Structured (2), Unstructured (4). |
| <b>Replace Fill Value With Nan</b><br>( <i>ReplaceFillValueWithNan</i> ) | If on, any float or double variable read that has a _FillValue attribute will have that fill value replaced with a not-a-number (NaN) value. The advantage of setting these to NaN values is that, if implemented properly by the system and careful math operations are used, they can implicitly be ignored by calculations like finding the range of the values. That said, this option should be used with caution as VTK does not fully support NaN values and therefore odd calculations may occur. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Spherical Coordinates</b><br>( <i>SphericalCoordinates</i> )          | If on, then data with latitude/longitude dimensions will be read in as curvilinear data shaped like spherical coordinates. If false, then the data will always be read in Cartesian coordinates.  | 1                | Only the values 0 and 1 are accepted.   |

|   |   |   |  |
|---|---|---|--|
| <b>Vertical Bias</b><br>( <i>VerticalBias</i> )   | The bias of the vertical component of spherical coordinates. It is common to write the vertical component with respect to something other than the center of the sphere (for example, the surface). In this case, it might be necessary to scale and/or bias the vertical height. The height will become height*scale + bias. Keep in mind that if the positive attribute of the vertical dimension is down, then the height is negated. The scaling will be adjusted if it results in invalid (negative) vertical values.  | 0 |  |
| <b>Vertical Scale</b><br>( <i>VerticalScale</i> ) | The scale of the vertical component of spherical coordinates. It is common to write the vertical component with respect to something other than the center of the sphere (for example, the surface). In this case, it might be necessary to scale and/or bias the vertical height. The height will become height*scale + bias. Keep in mind that if the positive attribute of the vertical dimension is down, then the height is negated. The scaling will be adjusted if it results in invalid (negative) vertical values. | 1 |  |

## Nrrd Reader

Read raw image files with Nrrd meta data.

The Nrrd reader reads raw image data much like the Raw Image Reader except that it will also read metadata information in the Nrrd format.

This means that the reader will automatically set information like file dimensions.

There are several limitations on what type of nrrd files we can read. This reader only supports nrrd files in raw format. Other encodings like ascii and hex will result in errors. When reading in detached headers, this only supports reading one file that is detached.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>Data VOI</b><br>( <i>DataVOI</i> )   | The data volume of interest (VOI). The VOI is a sub-extent of the data that you want loaded. Setting a VOI is useful when reading from a large data set and you are only interested in a small portion of the data. If left containing all 0's, then the reader will load in the entire data set. | 0 0 0 0 0 0      |              |
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to read (or the meta data file that will point to the actual file).  |                  |              |

## OpenFOAMReader

Read OpenFOAM data files, producing multi-block dataset.

The OpenFOAM reader reads OpenFOAM data files and outputs multi-block datasets. Mesh information and time dependent data are supported.

The OpenFOAM format is described fully at <http://www.openfoam.com/docs/user/basic-file-format.php>

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Add dimensional units to array names</b><br>( <i>AddDimensionsToArrayNames</i> )         | Read dimensional units from field data and add them to array names as human-readable string. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Cache mesh</b><br>( <i>CacheMesh</i> )   | Cache the OpenFOAM mesh between GUI selection changes.                                       | 1                | Only the values 0 and 1 are accepted.  |
| <b>Case Type</b><br>( <i>CaseType</i> )   | The property indicates whether decomposed mesh or reconstructed mesh should be read          | 1                | The value must be one of the following: Decomposed Case (0), Reconstructed Case (1). |
| <b>Cell Arrays</b><br>( <i>CellArrays</i> )   |  |                  | The list of array names is provided by the reader.                                   |
| <b>Create cell-to-point filtered data</b><br>( <i>CreateCellToPoint</i> )                   | Create point data from cell data. Beware: the filter does not do inverse distance weighting. | 1                | Only the values 0 and 1 are accepted.  |
| <b>Decompose polyhedra</b><br>( <i>DecomposePolyhedra</i> )                                 | Decompose polyhedra into tetrahedra and pyramids.  | 1                | Only the values 0 and 1 are accepted.  |
| <b>File Name</b><br>( <i>FileName</i> )   | This property specifies the file name for the reader.  |                  |  |
| <b>Lagrangian Arrays</b><br>( <i>LagrangianArrays</i> )                                     |  |                  | The list of array names is provided by the reader.                                   |
| <b>List timesteps according to controlDict</b><br>( <i>ListTimeStepsByControlDict</i> )     | List time directories listed according to the settings in controlDict.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Mesh Regions</b><br>( <i>MeshRegions</i> )   |  |                  | The list of array names is provided by the reader.                                   |
| <b>Point Arrays</b><br>( <i>PointArrays</i> )   |  |                  | The list of array names is provided by the reader.                                   |
| <b>Lagrangian positions are in OF 1.3 binary format</b><br>( <i>PositionsIsIn13Format</i> ) | Set if Lagrangian positions files are in OpenFOAM 1.3 binary format.                         | 0                | Only the values 0 and 1 are accepted.  |
| <b>Read zones</b><br>( <i>ReadZones</i> )   | Read point/face/cell-Zones?  | 0                | Only the values 0 and 1 are accepted.  |

## PDB Reader

Read PDB molecule files.

The PDB reader reads files in the format used by the Protein Data Bank (an archive of experimentally determined three-dimensional structures of biological macromolecules). The expected file extension is .pdb. The output datasets are polygonal.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the PDB reader. |                  |              |

## PLOT3D Reader

Read ASCII or binary PLOT3D files.

PLOT3D is a plotting package developed at NASA. The PLOT3D reader can read both ASCII and binary PLOT3D files. The default file extension for the geometry files is .xyz, and the default file extension for the solution files is .q. The output of this reader is a multi-block dataset containing curvilinear (structured grid) datasets.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Q File Name</b><br>( <i>QFileName</i> )                           | The list of .q (solution) files for the PLOT3D reader. This can be more than 1. If more than 1 file is specified, the reader will switch to file-series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |   |
| <b>Binary File</b><br>( <i>BinaryFile</i> )                          | This property indicates whether the data files are binary.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Byte Order</b><br>( <i>ByteOrder</i> )                            | This property specifies the byte order of any binary file(s).  | 0                | The value must be one of the following: BigEndian (0), LittleEndian (1).  |
| <b>File Name</b><br>( <i>FileName</i> )                              | This property specifies the .xyz (geometry) file name for the PLOT3D reader.   |                  |   |
| <b>Force Read</b><br>( <i>ForceRead</i> )                            | This property indicates whether to attempt to read the file even if it appears corrupt   | 0                | Only the values 0 and 1 are accepted.   |
| <b>Functions</b><br>( <i>Functions</i> )                             | This property lists which functions (point-centered arrays) to compute.  |                  | The value must be one of the following: Scalar - Pressure (110), Scalar - Temperature (120), Scalar - Enthalpy (130), Scalar - Internal energy (140), Scalar - Kinetic energy (144), Scalar - Velocity magnitude (153), Scalar - Entropy (170), Scalar - Swirl. (184), Vector - Velocity (200), Vector - Vorticity (201), Vector - Pressure gradient (210). |
| <b>Has Byte Count</b><br>( <i>HasByteCount</i> )                     | This property indicates whether this file uses Fortran leading-trailing byte counts.   | 0                | Only the values 0 and 1 are accepted.   |
| <b>I Blanking</b><br>( <i>IBlanking</i> )                            | This property indicates whether this dataset has visibility information.   | 0                | Only the values 0 and 1 are accepted.   |
| <b>Multi Grid</b><br>( <i>MultiGrid</i> )                            | This property indicates whether these files have multiple-block information.   | 0                | Only the values 0 and 1 are accepted.   |
| <b>Two Dimensional Geometry</b><br>( <i>TwoDimensionalGeometry</i> ) | This property indicates whether the data is two- or three-dimensional.   | 0                | Only the values 0 and 1 are accepted.   |

## PLY Reader

Read files stored in Stanford University's PLY polygonal file format.

The PLY reader reads files stored in the PLY polygonal file format developed at Stanford University. The PLY files that ParaView can read must have the elements "vertex" and "face" defined. The "vertex" elements must have the properties "x", "y", and "z". The "face" elements must have the property "vertex\_indices" defined. The default file extension for this reader is .ply.

| Property | Description | Default Value(s) | Restrictions |
|----------|-------------|------------------|--------------|
|----------|-------------|------------------|--------------|

|   |   |  |  |
|---|---|--|--|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the PLY reader. |  |  |
|---|---|--|--|

## PNG Series Reader

Read a PNG file into an image data.

The PNG reader reads PNG (Portable Network Graphics) files, and the output is a uniform rectilinear (image/volume) dataset. The default file extension is .png.

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## POP Reader

Read data files from the Parallel Ocean Program (POP).

The POP reader reads data files from the Parallel Ocean Program (POP). The expected file extension is .pop. The output of this reader is a structured grid dataset.

| Property  | Description  | Default Value(s)  | Restrictions |
|---|--|---|--------------|
| <b>Clip Extent</b><br>( <i>ClipExtent</i> )                     | This property indicates the minimum and maximum extent in each dimension of the dataset (X, Y, and Z) to read. | -2147483647 2147483647 -2147483647<br>2147483647 -2147483647 2147483647 |              |
| <b>File Name</b><br>( <i>FileName</i> )                         | This property specifies the file name for the POP Ocean file reader.   |   |              |
| <b>Number Of Ghost Levels</b><br>( <i>NumberOfGhostLevels</i> ) | This property indicates the number of ghost levels to generate.  | 1   |              |
| <b>Radius</b><br>( <i>Radius</i> )                              | This property specifies the radius of the dataset.   | 60000   |              |

## PVD Reader

Load a dataset stored in ParaView's PVD file format.

The PVD reader reads data stored in ParaView's PVD file format. The .pvd file is essentially a header file that collects together other data files stored in VTK's XML-based file format.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the PVD reader. |                  |              |

## Particles Reader

Read particle data.

vtkParticleReader reads either a binary or a text file of particles.

Each particle can have associated with it an optional scalar value. So

the format is: x, y, z, scalar (all floats or doubles). The text file

can consist of a comma delimited set of values. In most cases

vtkParticleReader can automatically determine whether the file is text

or binary. The data can be either float or double. Progress updates

are provided. With respect to binary files, random access into the file

to read pieces is supported.

| Property                                  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )   | The list of files to be read by the reader.   |                  |  |
| <b>Byte Order</b><br>( <i>ByteOrder</i> ) | Set/Get the byte ordering of the file you are trying to read in.  | 0                | The value must be one of the following: BigEndian (0), LittleEndian (1).   |
| <b>Data Type</b><br>( <i>DataType</i> )   | Get/Set the data type. The options are: - Float (10) (default) single precision floating point. - Double (11) double precision floating point.  | 10               | The value must be one of the following: Float (10), Double (11).           |
| <b>File Type</b><br>( <i>FileType</i> )   | Set the file type. The options are: - Unknown (0) (default) the reader will attempt to determine the file type. If this fails then you should set the file type yourself. - Text (1) the file type is text. - Binary (2) the file type is binary. | 0                | The value must be one of the following: Unknown (0), Text (1), Binary (2). |
| <b>Has Scalar</b><br>( <i>HasScalar</i> ) | Default: 1. If 1 then each particle has a value associated with it.   | 1                | Only the values 0 and 1 are accepted.                                      |

## Partitioned Legacy VTK Reader

Read files stored in VTK partitioned legacy format.

The Partitioned Legacy VTK reader loads files stored in VTK's partitioned legacy file format (before VTK 4.2, although still supported). The expected file extension is .pvtk. The type of the dataset may be structured grid, uniform rectilinear grid (image/volume), non-uniform rectilinear grid, unstructured grid, or polygonal.

| Property                                | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the Partitioned Legacy VTK reader. |                  |              |



## Phasta Reader

Read the parallel Phasta meta-file and the underlying Phasta files.

This Phasta reader reads files stored in the Phasta (a CFD package) format. The expected file extension is .pht. The output of this reader is a multipiece data set.

| Property                                | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the Phasta reader. |                  |              |

## Restarted Sim Exodus Reader

Read collections of Exodus output files from simulations that were restarted.

When a simulation that outputs exodus files is restarted, typically you get a new set of output files. When you read them in your visualization, you often want to string these file sets together as if it was one continuous dump of files. This reader allows you to specify a metadata file that will implicitly string the files together.

| Property  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                           | This points to a special metadata file that lists the output files for each restart.  |                  |  |
| <b>Use Meta File</b><br>( <i>UseMetaFile</i> )                    | This hidden property must always be set to 1 for this proxy to work.  | 1                | Only the values 0 and 1 are accepted.              |
| <b>Animate Vibrations</b><br>( <i>AnimateVibrations</i> )         | If this flag is on and HasModeShapes is also on, then this reader will report a continuous time range [0,1] and animate the displacements in a periodic sinusoid. If this flag is off and HasModeShapes is on, this reader ignores time. This flag has no effect if HasModeShapes is off.   | 1                | Only the values 0 and 1 are accepted.              |
| <b>Apply Displacements</b><br>( <i>ApplyDisplacements</i> )       | Geometric locations can include displacements. When this option is on, the nodal positions are 'displaced' by the standard exodus displacement vector. If displacements are turned 'off', the user can explicitly add them by applying a warp filter.   | 1                | Only the values 0 and 1 are accepted.              |
| <b>Displacement Magnitude</b><br>( <i>DisplacementMagnitude</i> ) | When displacements are being applied, they are scaled by this amount. Set to 1 for no scaling.  | 1                |  |
| <b>Edge Blocks</b><br>( <i>EdgeBlocks</i> )                       | Use this property to select which blocks of edges should be loaded. All edges in the same block will have the same interpolant (linear, quadratic, etc.) and have the same variables defined over them.   |                  | The list of array names is provided by the reader. |
| <b>Edge Map Array Status</b><br>( <i>EdgeMapArrayStatus</i> )     | Specify which maps defined over edges in the mesh should be loaded. An edge map is a renumbering of all the edges (across all the edge blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files.   |                  | The list of array names is provided by the reader. |
| <b>Edge Set Array Status</b><br>( <i>EdgeSetArrayStatus</i> )     | An Exodus II file may define subsets of all the edges (across all edge blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per edge. The accompanying EdgeSetResultArrayStatus property specifies which variables defined over those sets should be loaded. |                  | The list of array names is provided by the reader. |

|  |   |   |  |
|--|---|---|--|
| <b>Edge Set Result Array Status</b><br>( <i>EdgeSetResultArrayStatus</i> )       | An Exodus II file may define subsets of all the edges (across all edge blocks) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per edge. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the EdgeSetArrayStatus property in order for the variable to appear.                          |   | The list of array names is provided by the reader. |
| <b>Edge Variables</b><br>( <i>EdgeVariables</i> )                                | Use this property to select which variables defined over edges should be loaded. Edges are oriented boundaries of one or more higher-dimensional cells also present in the mesh. Edge variables specify a single value per entry in some edge block.  |   | The list of array names is provided by the reader. |
| <b>Element Blocks</b><br>( <i>ElementBlocks</i> )                                | Use this property to select which blocks of elements (cells) should be loaded. All elements (cells) in the same block will have the same interpolant (tetrahedral, triangular, trilinear hexahedral, etc.) and have the same variables defined over them.   |   |  |
| <b>Element Map Array Status</b><br>( <i>ElementMapArrayStatus</i> )              | Specify which maps defined over elements (cells) in the mesh should be loaded. An element map is a renumbering of all the elements (across all the element blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files. These are sometimes called new-style maps because they are specified differently than in older Exodus files, which only allowed a single element map. |   | The list of array names is provided by the reader. |
| <b>Element Set Array Status</b><br>( <i>ElementSetArrayStatus</i> )              | An Exodus II file may define subsets of all the elements (across all element blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as feature classifiers, may then be defined over these sets by specifying a single number per element. The accompanying ElementSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                 |   | The list of array names is provided by the reader. |
| <b>Element Set Result Array Status</b><br>( <i>ElementSetResultArrayStatus</i> ) | An Exodus II file may define subsets of all the elements (across all element blocks) in a file as sets in their own right. Variables, such as feature classifiers, may then be defined over these sets by specifying a single number per element. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the ElementSetArrayStatus property in order for the variable to appear.              |   | The list of array names is provided by the reader. |
| <b>Element Variables</b><br>( <i>ElementVariables</i> )                          | Use this property to select which variables defined over elements (cells) should be loaded. Element variables specify a single value per entry in some element block.   |   | The list of array names is provided by the reader. |
| <b>Exodus Model Metadata</b><br>( <i>ExodusModelMetadata</i> )                   | Should metadata for the mesh be stored in the field data of the reader's output?  | 0 | Only the values 0 and 1 are accepted.              |
| <b>Face Blocks</b><br>( <i>FaceBlocks</i> )                                      | Use this property to select which blocks of faces should be loaded. All faces in the same block will have the same interpolant (triangular, bilinear quadrilateral, etc.) and have the same variables defined over them.  |   | The list of array names is provided by the reader. |
| <b>Face Map Array Status</b><br>( <i>FaceMapArrayStatus</i> )                    | Specify which maps defined over faces in the mesh should be loaded. A face map is a renumbering of all the faces (across all the face blocks of a file) from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files.  |   | The list of array names is provided by the reader. |
| <b>Face Set Array Status</b><br>( <i>FaceSetArrayStatus</i> )                    | An Exodus II file may define subsets of all the faces (across all face blocks) in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per face. The accompanying FaceSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                             |   | The list of array names is provided by the reader. |

|  |  |   |  |
|--|--|---|--|
| <b>Face Set Result Array Status</b><br>( <i>FaceSetResultArrayStatus</i> )         | An Exodus II file may define subsets of all the faces (across all face blocks) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per face. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the FaceSetArrayStatus property in order for the variable to appear. |   | The list of array names is provided by the reader. |
| <b>Face Variables</b><br>( <i>FaceVariables</i> )                                  | Use this property to select which variables defined over faces should be loaded. Faces are oriented boundaries of one or more higher-dimensional cells also present in the mesh. Face variables specify a single value per entry in some face block.   |   | The list of array names is provided by the reader. |
| <b>Generate Global Element Id Array</b><br>( <i>GenerateGlobalElementIdArray</i> ) | Should the reader retrieve the first new-style element map, or if that is not present, the solitary old-style element map (which always exists but may be procedurally generated if it is not stored with the file)?   | 1 | Only the values 0 and 1 are accepted.              |
| <b>Generate Global Node Id Array</b><br>( <i>GenerateGlobalNodeIdArray</i> )       | Should the reader retrieve the first new-style node map, or if that is not present, the solitary old-style node map (which always exists but may be procedurally generated if it is not stored with the file)?   | 1 | Only the values 0 and 1 are accepted.              |
| <b>Generate Object Id Cell Array</b><br>( <i>GenerateObjectIdCellArray</i> )       | Specifies whether a variable named "ObjectId" should be created for each entry of each block and set being loaded and assigned the identifying number of the block or set, so that given a cell one can readily identify the block or set it came from. This is useful for coloring a mesh by block ID or set ID.  | 1 | Only the values 0 and 1 are accepted.              |
| <b>Global Variables</b><br>( <i>GlobalVariables</i> )                              | Use this property to select which global (i.e., per-mesh) variables should be loaded. Global variables are stored as arrays in the field data of the reader's output.  |   | The list of array names is provided by the reader. |
| <b>Has Mode Shapes</b><br>( <i>HasModeShapes</i> )                                 | Some simulations overload the Exodus time steps to represent mode shapes. In this case, it does not make sense to iterate over the "time steps", because they are not meant to be played in order. Rather, each represents the vibration at a different "mode." Setting the value of this property to 1 changes the semantics of the reader to not report the time steps to downstream filters.                                  | 0 | Only the values 0 and 1 are accepted.              |
| <b>Mode Shape</b><br>( <i>ModeShape</i> )  | Specify which mode shape to animate when HasModeShapes is on.  | 1 |  |
| <b>Node Map Array Status</b><br>( <i>NodeMapArrayStatus</i> )                      | Specify which maps defined over nodes (points) in the mesh should be loaded. A node map is a renumbering of all the points in a file from their offset in the file to some arbitrary integers, usually specifying a unique global ID across a set of files. These are sometimes called new-style maps because they are specified differently than in older Exodus files, which only allowed a single node map.                   |   | The list of array names is provided by the reader. |
| <b>Node Set Array Status</b><br>( <i>NodeSetArrayStatus</i> )                      | An Exodus II file may define subsets of all the nodes (points) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per node. This property specifies which of those sets should be loaded. The accompanying NodeSetResultArrayStatus property specifies which variables defined over those sets should be loaded.                    |   | The list of array names is provided by the reader. |
| <b>Node Set Result Array Status</b><br>( <i>NodeSetResultArrayStatus</i> )         | An Exodus II file may define subsets of all the nodes (points) in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per node. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the NodeSetArrayStatus property in order for the variable to appear.                 |   | The list of array names is provided by the reader. |
| <b>Point Variables</b><br>( <i>PointVariables</i> )                                | Use this property to select which nodal (i.e., per-point) variables should be loaded.  |   | The list of array names is provided by the reader. |

|  |   |  |  |
|--|---|--|--|
| <b>Side Set Array Status</b><br>( <i>SideSetArrayStatus</i> )              | An Exodus II file may define subsets of all the of all the elements in a file as sets in their own right. This property specifies which of those sets should be loaded. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per side. For example, a hexahedron has 18 sides: 6 faces and 12 edges. Any of these sides may be individually called out in a set and assigned a result value. The accompanying SideSetResultArrayStatus property specifies which variables defined over those sets should be loaded.    |  | The list of array names is provided by the reader. |
| <b>Side Set Result Array Status</b><br>( <i>SideSetResultArrayStatus</i> ) | An Exodus II file may define subsets of all the of all the elements in a file as sets in their own right. Variables, such as boundary conditions, may then be defined over these sets by specifying a single number per side. For example, a hexahedron has 18 sides: 6 faces and 12 edges. Any of these sides may be individually called out in a set and assigned a result value. This property specifies which of those variables should be loaded. The corresponding set must also be loaded using the SideSetArrayStatus property in order for the variable to appear. |  | The list of array names is provided by the reader. |
| <b>XML File Name</b><br>( <i>XMLFileName</i> )                             | This property specifies the name of an XML file containing part and material descriptions and a mapping to element blocks.  |  |  |

## Restarted Sim Spy Plot Reader

Read collections of SPCTH files from simulations that were restarted.

When a CTH simulation is restarted, typically you get a new set of output files. When you read them in your visualization, you often want to string these file sets together as if it was one continuous dump of files. This reader allows you to specify a metadata file that will implicitly string the files together.

| Property  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                     | This points to a special metadata file that lists the output files for each restart.                            |                  |  |
| <b>Use Meta File</b><br>( <i>UseMetaFile</i> )                              | This hidden property must always be set to 1 for this proxy to work.  | 1                | Only the values 0 and 1 are accepted.              |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )                            | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Distribute Files</b><br>( <i>DistributeFiles</i> )                       | In parallel mode, if this property is set to 1, the reader will distribute files or blocks.                     | 1                | Only the values 0 and 1 are accepted.              |
| <b>Down Convert Volume Fraction</b><br>( <i>DownConvertVolumeFraction</i> ) | If this property is set to 0, the type of the volume fraction is float; is set to 1, the type is unsigned char. | 1                | Only the values 0 and 1 are accepted.              |
| <b>Generate Active Block Array</b><br>( <i>GenerateActiveBlockArray</i> )   | If this property is set to 1, a cell array will be generated that stores the active status of a block.          | 0                | Only the values 0 and 1 are accepted.              |
| <b>Generate Block Id Array</b><br>( <i>GenerateBlockIdArray</i> )           | If this property is set to 1, a cell array will be generated that stores a unique blockId for each block.       | 0                | Only the values 0 and 1 are accepted.              |
| <b>Generate Level Array</b><br>( <i>GenerateLevelArray</i> )                | If this property is set to 1, a cell array will be generated that stores the level of each block.               | 0                | Only the values 0 and 1 are accepted.              |

## SESAME Reader

Read SESAME data files, producing rectilinear grids.

The SESAME reader reads SESAME data files, and outputs rectilinear grids. The expected file extension is .sesame.

| Property                                | Description  | Default Value(s) | Restrictions                             |
|---|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the SESAME reader. |                  |  |
| <b>Table Id</b><br>( <i>TableId</i> )   | This property indicates which table to read.                 | -1               | The value must be one of the following:. |

## SLAC Data Reader

A reader for a data format used by Omega3p, Tau3p, and several other tools used at the Stanford Linear Accelerator Center (SLAC).

The underlying format uses netCDF to store arrays, but also imposes several conventions to form an unstructured grid of elements.

| Property   | Description  | Default Value(s) | Restrictions                          |
|--|--|------------------|---------------------------------------|
| <b>Mesh File Name</b><br>( <i>MeshFileName</i> )               | The name of the mesh file to load.   |                  |                                       |
| <b>Mode File Name</b><br>( <i>ModeFileName</i> )               | The name of the mode files to load. The points in the mode file should be the same as the mesh file. |                  |                                       |
| <b>Read External Surface</b><br>( <i>ReadExternalSurface</i> ) | If on, read the external surfaces of the data set.   | 1                | Only the values 0 and 1 are accepted. |
| <b>Read Internal Volume</b><br>( <i>ReadInternalVolume</i> )   | If on, read the internal volume of the data set.   | 0                | Only the values 0 and 1 are accepted. |
| <b>Read Midpoints</b><br>( <i>ReadMidpoints</i> )              | If on, reads midpoint information for external surfaces and builds quadratic surface triangles.      | 0                | Only the values 0 and 1 are accepted. |

## SLAC Particle Data Reader

The SLAC Particle data reader.

| Property                                | Description                                  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | A list of files to be read in a time series. |                  |              |

## STL Reader

Read ASCII or binary stereo lithography (STL) files.

The STL reader reads ASCII or binary stereo lithography (STL) files. The expected file extension is .stl. The output of this reader is a polygonal dataset. This reader also supports file series.

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## Spy Plot History Reader

Read a spth history file where each row translates into a single time step and the columns are points, materials and properties.

| Property  | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>Comment Character</b><br>( <i>CommentCharacter</i> ) | This property lists the characters that is used as the first character on comment lines | %                |              |
| <b>Delimiter</b><br>( <i>Delimiter</i> )                | Character that is used as the delimiter.  | ,                |              |
| <b>File Name</b><br>( <i>FileName</i> )                 | This property specifies the file name for the VRML reader.                              |                  |              |

## Spy Plot Reader

Read files in the SPCTH Spy Plot file format.

The Spy Plot reader loads an ASCII meta-file called the "case" file (extension .spcth). The case file lists all the binary files containing the dataset. This reader produces hierarchical datasets.

| Property  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )                            | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Distribute Files</b><br>( <i>DistributeFiles</i> )                       | In parallel mode, if this property is set to 1, the reader will distribute files or blocks.                     | 1                | Only the values 0 and 1 are accepted.              |
| <b>Down Convert Volume Fraction</b><br>( <i>DownConvertVolumeFraction</i> ) | If this property is set to 0, the type of the volume fraction is float; is set to 1, the type is unsigned char. | 1                | Only the values 0 and 1 are accepted.              |
| <b>File Name</b><br>( <i>FileName</i> )                                     | This property specifies the file name for the Spy Plot reader.  |                  |  |
| <b>Generate Active Block Array</b><br>( <i>GenerateActiveBlockArray</i> )   | If this property is set to 1, a cell array will be generated that stores the active status of a block.          | 0                | Only the values 0 and 1 are accepted.              |
| <b>Generate Block Id Array</b><br>( <i>GenerateBlockIdArray</i> )           | If this property is set to 1, a cell array will be generated that stores a unique blockId for each block.       | 0                | Only the values 0 and 1 are accepted.              |
| <b>Generate Level Array</b><br>( <i>GenerateLevelArray</i> )                | If this property is set to 1, a cell array will be generated that stores the level of each block.               | 0                | Only the values 0 and 1 are accepted.              |

|  |   |   |                                       |
|--|---|---|---------------------------------------|
| <b>Generate Tracer Array</b><br>( <i>GenerateTracerArray</i> ) | If this property is set to 1, a cell array will be generated that stores the coordinates of any tracers.  | 0 | Only the values 0 and 1 are accepted. |
| <b>Merge XYZComponents</b><br>( <i>MergeXYZComponents</i> )    | If this property is set to 1, cell arrays named (for example) X velocity, Y velocity and Z velocity will be combined into a single vector array named velocity. | 1 | Only the values 0 and 1 are accepted. |

## TIFF Reader

Read a TIFF file into an image data.

The TIFF reader reads TIFF (Tagged Image File Format) files, and the output is a uniform rectilinear (image/volume) dataset. The default file extension is .tiff.

| Property                                | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the TIFF reader. |                  |              |

## TIFF Series Reader

Read a series of TIFF files into an time sequence of image datas.

The TIFF series reader reads TIFF files. The output is a time sequence of uniform rectilinear (image/volume) dataset. The default file extension is .tif or .tiff.

| Property                                  | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Names</b><br>( <i>FileNames</i> ) | The list of files to be read by the reader. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## Tecplot Reader

Read files in the Tecplot ASCII file format.

The Tecplot reader extracts multiple zones (blocks) of data from a Tecplot ASCII file, in which a zone is stored in either point packing mode (i.e., tuple-based, with only point data supported) or block packing mode (i.e., component-based, with point data and cell data supported). The output of the reader is a vtkMultiBlockDataset, of which each block is either a vtkStructuredGrid or a vtkUnstructuredGrid. This supports reading a file series.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Names</b><br>( <i>FileNames</i> )        | The list of files to be read by the reader.   |                  |  |
| <b>Data Arrays</b><br>( <i>dataArrayStatus</i> ) | This property lists what dataset attribute arrays (point data and cell data) to read. |                  | The list of array names is provided by the reader. |

## VPIC Reader

Read distributed VPIC files into an ImageData.

VPIC is a 3D kinetic plasma particle-in-cell simulation. The input file (.vpc) opened by the VPIC reader is an ASCII description of the data files which are written one file per processor, per category and per time step. These are arranged in subdirectories per category (field data and hydrology data) and then in time step subdirectories.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | ASCII .vpc file describes locations of data files, grid sizes, time step sizes and type and order of data written to the files.   |                  |  |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | Variables written to the data files are described in the .vpc file and are presented for selection. Only selected variables are loaded for a time step.   |                  | The list of array names is provided by the reader. |
| <b>Stride</b><br>( <i>SetStride</i> )              | VPIC data may be very large and not all is needed for effective visualization. Setting the stride selects every nth data item within the files for display.   | 1 1 1            | The value must be greater than or equal to 1.      |
| <b>X Extent</b><br>( <i>XExtent</i> )              | VPIC data is written one file per simulation processor. This coarse map of files is used in partitioning files between visualizing processors so that each ParaView processor has its own set of files to display. Ghost cell overlap is handled within the reader. To limit the View of VPIC information the extent in the X dimension of "files" can be specified. Only the files selected will be displayed and they will be partitioned between the visualizing processors, allowing a higher resolution over a smaller area. | -1 -1            |  |
| <b>Y Extent</b><br>( <i>YExtent</i> )              | VPIC data is written one file per simulation processor. This coarse map of files is used in partitioning files between visualizing processors so that each ParaView processor has its own set of files to display. Ghost cell overlap is handled within the reader. To limit the View of VPIC information the extent in the Y dimension of "files" can be specified. Only the files selected will be displayed and they will be partitioned between the visualizing processors, allowing a higher resolution over a smaller area. | -1 -1            |  |
| <b>Z Extent</b><br>( <i>ZExtent</i> )              | VPIC data is written one file per simulation processor. This coarse map of files is used in partitioning files between visualizing processors so that each ParaView processor has its own set of files to display. Ghost cell overlap is handled within the reader. To limit the View of VPIC information the extent in the Z dimension of "files" can be specified. Only the files selected will be displayed and they will be partitioned between the visualizing processors, allowing a higher resolution over a smaller area. | -1 -1            |  |

## VRML Reader

Load the geometry from a VRML 2.0 file.

The VRML reader loads only the geometry from a VRML (Virtual Reality Modeling Language) 2.0 file. The expected file extension is .wrl. The output of this reader is a polygonal dataset.

| Property                                | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the VRML reader. |                  |              |



## VisItANALYZEReader

Format that stores fMRI data of brain scans

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .img and .hdr.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItANSYSReader

ASCII file format used for input to the ANSYS suite

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .inp.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItAUXFileReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .aux.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItBOVReader

BOV (Brick of Values) generates a rectilinear grid of values

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .bov.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItCEAucdReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions .ucd, and .inp

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItCMATReader

ASCII output of CMAT Fortran simulation.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .cmat

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItCTRLReader

Reads control files for TB-LMTO

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .ctrl

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItChomboReader

Reads files from the Chombo package. Loads block-structured adaptively refined rectangular grids

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .h5 and .hdf5.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItClawReader

A material-ordered format very similar to IOUT=1.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .claw

| Property                                  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>Cell Arrays</b><br>(CellArrayStatus)   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>(FileName)            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>(MaterialStatus)      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>(MeshStatus)             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>(PointArrayStatus) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItCurve2DReader

Reads ASCII 2D Curve Format. Which is a single x,y values per line to generate a 2D curve

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .curve, .ultra, .ult, and .u

| Property                                  | Description   | Default Value(s) | Restrictions                                       |
|---|---|------------------|--|
| <b>Cell Arrays</b><br>(CellArrayStatus)   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>(FileName)            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>(MaterialStatus)      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>(MeshStatus)             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>(PointArrayStatus) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItDDCMDReader

Reads ddcMD particle and atom files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .ddcmd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItDyna3DReader

Reads Dyna3D finite element files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .dyn

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItEnzoReader

Reads enzo cosmological simulation AMR files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .boundary, and .hierarchy

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItExtrudedVolReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .exvol.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItFLASHReader

ASC / Alliances Center for Astrophysical Thermonuclear Flash Reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .flash, and .f5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItFluentReader

Fluent CFD simulation result reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .cas

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |



## VisItGGCMReader

Reads Geospace General Circulation Model files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .3df, and .mer

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItGTCReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItGULPReader

GULP trajectory file reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .trg

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItGadgetReader

Gadget 2 cosmological N-body/SPH simulation code reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .gadget

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItH5NimrodReader

NIMROD HDF5 data model reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5nimrod

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItImageReader

Multiple Image format reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .pnm, .ppm, .sdt, .spr and .imgvol

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItLAMMPSDumpReader

LAMMPS Molecular Dynamics Simulator dump reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .dump

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItLAMMPSStructureReader

LAMMPS Molecular Dynamics Simulator reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .eam, .meam, .rigid, and .lammgs

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItLinesReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .lines.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItM3DC1Reader

M3DC1 fusion code reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItM3DReader

M3D fusion code reader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .m3d

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItMM5Reader

Reads PSU/NCAR mesoscale model simulation files.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is mm5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItMirandaReader

Readers the raw output format of LLNL's Miranda simulation.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .mir, and .raw

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNASTRANReader

Reads NASTRAN Bulk ASCII unstructured meshes

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nas, and .f06

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNek5000Reader

Reads simulation results from Nek5000 which is a DNS/LES computational fluid dynamics solver

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nek3, .nek2d, .nek5d, .nek5000, .nek

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFADAPTReader

Reads NetCDF ADAPT files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |



## VisItNetCDFCCSM\_MTSDReader

Reads CCSM climate data files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFCCSM\_STSDReader

Reads CCSM climate data files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFFVCOMParticleReader

Reads FVCOM primitive equations ocean model particle files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFFVCOM\_MTMDReader

Reads FVCOM primitive equations ocean models

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDDFFVCOM\_MTSDReader

Reads FVCOM primitive equations ocean models

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDDFFVCOM\_STSDReader

Reads FVCOM primitive equations ocean models

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFLODIParticleReader

Reads NetCDF LODI particle files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItNetCDFLODIReader

Reads NetCDF LODI files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .nc, .cdf, .elev, and .ncd

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItOVERFLOWReader

Reads Binary Fortran output files containing overlaid curvilinear meshes files.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .dat, and .save

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItOpenFOAMReader

Reads simulation results from OpenFOAM (Open Field Operation and Manipulation) CFD Toolbox

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .controlDict

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPATRANReader

Reads PATRAN neutral files, which are ASCII format files that contain unstructured geometry and simulation results

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .neu

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPFLOTTRANReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPLOT2DReader

Reads structured simulation data from Plot2D.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .p2d

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPLOT3DReader

Reads structured simulation data from Plot3D.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .q, .x, and .vp3d

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItParaDisReader

Reads Parallel Dislocation Simulator (ParaDiS) files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .prds, .data, .dat

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItParaDisTecplotReader

Reads Parallel Dislocation Simulator (ParaDiS) tecplot like files

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .fld, .field, .cyl, .cylinder, .dat

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |



## VisItPixieReader

Reads 2D and 3D simulation data saved in Pixie HDF5 format.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPlainTextReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .csv, .tsv, and .txt.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItPoint3DReader

Reads an ASCII file containing X,Y,Z, and data value columns.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .3D

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItProteinDataBankReader

Reads Protein Databank files contain molecular data.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .ent, and .pdb

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItRAWReader

Reads RAW mesh geometry format.

Allows mesh geometry to be shared with CAD programs such as Blender and POV-Ray

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .raw

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItSAMRAIReader

Reads Structured AMR data stored in HDF5 file formats.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .samrai

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItSARReader

Reads Specific Absorption Rate image volume files.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .sar

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItSASReader

Reads SAS files.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .sas, sasgeom, and .sasdata

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItSphericalReader

Reads ASCII coupled hydrodynamical and gravitational simulations

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .spherical, and .sv

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItTFTReader

Reads ASCII TFT Fortran simulation results.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .dat, and .tft

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItTSurfReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .ts\_deg83.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItTecplotBinaryReader

Reads a subset of the Tecplot binary file format

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .plt

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItTecplotReader

Reads a subset of the ASCII Tecplot file format

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .tp, and .tec

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItTetradReader

Reads tetrahedral meshes and variables stored in HDF5 file format.

The default file extensions are .hdf5, and .h5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItUNICReader

Reads UNIC neutronics code result files.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .h5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItVASPCHGCarReader

Reads VASP simulation result files.

VAMP/VASP is a package for performing ab-initio quantum-mechanical molecular dynamics (MD) using pseudopotentials and a plane wave basis set.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is CHG\*

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItVASPOUTCARReader

Reads VASP simulation result files.

VAMP/VASP is a package for performing ab-initio quantum-mechanical molecular dynamics (MD) using pseudopotentials and a plane wave basis set.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is out\*



| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItVASPPOSCARReader

Reads VASP simulation result files.

VAMP/VASP is a package for performing ab-initio quantum-mechanical molecular dynamics (MD) using pseudopotentials and a plane wave basis set.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is POS\*

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItVelodyneReader

Reads Velodyne structural mechanics simulation result files.

Velodyne as an advanced structural mechanics code written by Corvid Technologies.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .vld, and .rst

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItVsReader

Reads VizSchema hdf5 files.

VizSchema is a specification for markup added to hdf5 files to describe the data contained within. VizSchema also allows users to specify run information and other miscellaneous data in a standardized manner.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extensions are .h5, and .vsh5

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItXYZReader

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .xyz.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## VisItXmdvReader

Reads XmdvTool ASCII files.

XmdvTool is a public-domain software package for the interactive visual exploration of multivariate data sets.

Note this reader is automatically generated from wrapping a third party reader. For more information on the reader see <https://wci.llnl.gov/codes/visit>

The default file extension is .okc

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.     |                  | The list of array names is provided by the reader. |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the reader        |                  |  |
| <b>Materials</b><br>( <i>MaterialStatus</i> )      | Select the Materials to be loaded from the dataset, if any. |                  | The list of array names is provided by the reader. |
| <b>Meshes</b><br>( <i>MeshStatus</i> )             | Select the meshes to be loaded from the dataset, if any.    |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.    |                  | The list of array names is provided by the reader. |

## Wavefront OBJ Reader

Read Wavefront .OBJ files to produce polygonal and line data.

The OBJ reader reads data stored in Wavefront .OBJ format.

The expected file extension is .obj, the datasets resulting from reading these files are polygons and lines.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the OBJ reader. |                  |              |

## WindBlade reader

Read WindBlade/Firetec simulation files possibly including wind turbines and topology files.

WindBlade/Firetec is a simulation dealing with the effects of wind on wind turbines or on the spread of fires. It produces three outputs - a StructuredGrid for the wind data fields, a StructuredGrid for the ground topology, and a PolyData for turning turbine blades. The input file (.wind) opened by the WindBlade reader is an ASCII description of the data files expected. Data is accumulated by the simulation processor and is written one file per time step. WindBlade can deal with topology if a flag is turned on and expects (x,y) data for the ground. It also can deal with turning wind turbines from other time step data files which gives polygon positions of segments of the blades and data for each segment.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>Filename</b><br>( <i>Filename</i> )             | ASCII .wind file describes locations of data files, grid sizes and variable deltas, time step sizes, whether topology is used, whether turbines are used, and type and order of data written to the files. |                  |  |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | Variables written to the data files are described in the .wind file and are presented for selection. Only selected variables are loaded for a time step.   |                  | The list of array names is provided by the reader. |

## XDMF Reader

Read XDMF (eXtensible Data Model and Format) files.

The XDMF reader reads files in XDMF format. The expected file extension is .xmf. Metadata is stored in the XDMF file using an XML format, and large attribute arrays are stored in a corresponding HDF5 file. The output may be unstructured grid, structured grid, or rectilinear grid. See <http://www.xdmf.org> for a description of the file format.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader.    |
| <b>File Name</b><br>( <i>FileName</i> )            | This property specifies the file name for the XDMF reader.  |                  |   |
| <b>Grids</b><br>( <i>GridStatus</i> )              | Controls which particular data sets to read from a file that contains many data sets inside a composite data set collection.                            |                  |   |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader.    |
| <b>Sets</b><br>( <i>SetStatus</i> )                | Select the sets to be loaded from the dataset, if any.  |                  | The list of array names is provided by the reader.    |
| <b>Stride</b><br>( <i>Stride</i> )                 | If loading structured data, this property indicate the number of indices per dimension (X, Y, or Z) to skip between each point included in this output. | 1 1 1            | The value must be greater than or equal to (1, 1, 1). |

## XML Hierarchical Box Data reader

Read a VTK XML-based data file containing a hierarchical dataset containing vtkUniformGrids.

The XML Hierarchical Box Data reader reads VTK's XML-based file format containing a vtkHierarchicalBoxDataSet. The expected file extensions is either .vthb or .vth.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | The list of files to be read by the reader. Each file is expected to be in the VTK XML polygonal dataset format. The standard extension is .vtp. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## XML Image Data Reader

Read serial VTK XML image data files.

The XML Image Data reader reads the VTK XML image data file format. The standard extension is .vti. This reader also supports file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the VTK XML image data format. The standard extension is .vti. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.   |                  | The list of array names is provided by the reader. |

## XML MultiBlock Data Reader

Read a VTK XML multi-block data file and the serial VTK XML data files to which it points.

The XML Multi-Block Data reader reads the VTK XML multi-block data file format. XML multi-block data files are meta-files that point to a list of serial VTK XML files. When reading in parallel, this reader will distribute sub-blocks among processors. The expected file extensions are .vtm and .vtmb.

| Property                                | Description   | Default Value(s) | Restrictions |
|---|---|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | The list of files to be read by the reader. Each file is expected to be in the VTK XML polygonal dataset format. The standard extension is .vtp. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |              |

## XML Partitioned Image Data Reader

Read the summary file and the associated VTK XML image data files.

The XML Partitioned Image Data reader reads the partitioned VTK image data file format. It reads the partitioned format's summary file and then the associated VTK XML image data files. The expected file extension is .pvti. This reader also supports file series.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the partitioned VTK XML image data format. The standard extension is .pvti. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## XML Partitioned Polydata Reader

Read the summary file and the associated VTK XML polydata files.

The XML Partitioned Polydata reader reads the partitioned VTK polydata file format. It reads the partitioned format's summary file and then the associated VTK XML polydata files. The expected file extension is .pvtp. This reader also supports file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the partitioned VTK XML polygonal dataset format. The standard extension is .pvtp. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.   |                  | The list of array names is provided by the reader. |

## XML Partitioned Rectilinear Grid Reader

Read the summary file and the associated VTK XML rectilinear grid data files.

The XML Partitioned Rectilinear Grid reader reads the partitioned VTK rectilinear grid file format. It reads the partitioned format's summary file and then the associated VTK XML rectilinear grid files. The expected file extension is .pvtr. This reader also supports file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the partitioned VTK XML rectilinear grid data format. The standard extension is .pvtr. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.   |                  | The list of array names is provided by the reader. |

## XML Partitioned Structured Grid Reader

Read the summary file and the associated VTK XML structured grid data files.

The XML Partitioned Structured Grid reader reads the partitioned VTK structured grid data file format. It reads the partitioned format's summary file and then the associated VTK XML structured grid data files. The expected file extension is .pvts. This reader also supports file series.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the partitioned VTK XML structured grid data format. The standard extension is .pvts. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## XML Partitioned Unstructured Grid Reader

Read the summary file and the associated VTK XML unstructured grid data files.

The XML Partitioned Unstructured Grid reader reads the partitioned VTK unstructured grid data file format. It reads the partitioned format's summary file and then the associated VTK XML unstructured grid data files. The expected file extension is .pvtu. This reader also supports file series.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the partitioned VTK XML unstructured grid data format. The standard extension is .pvtu. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## XML PolyData Reader

Read serial VTK XML polydata files.

The XML Polydata reader reads the VTK XML polydata file format. The standard extension is .vtp. This reader also supports file series.

| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the VTK XML polygonal dataset format. The standard extension is .vtp. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## XML Rectilinear Grid Reader

Read serial VTK XML rectilinear grid data files.

The XML Rectilinear Grid reader reads the VTK XML rectilinear grid data file format. The standard extension is .vtr. This reader also supports file series.



| Property   | Description   | Default Value(s) | Restrictions                                       |
|--|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the VTK XML rectilinear grid data format. The standard extension is .vtr. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.   |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.  |                  | The list of array names is provided by the reader. |

## XML Structured Grid Reader

Read serial VTK XML structured grid data files.

The XML Structured Grid reader reads the VTK XML structured grid data file format. The standard extension is .vts. This reader also supports file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the VTK XML structured grid data format. The standard extension is .vts. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.   |                  | The list of array names is provided by the reader. |

## XML Unstructured Grid Reader

Read serial VTK XML unstructured grid data files.

The XML Unstructured Grid reader reads the VTK XML unstructured grid data file format. The standard extension is .vtu. This reader also supports file series.

| Property   | Description  | Default Value(s) | Restrictions                                       |
|--|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )            | The list of files to be read by the reader. Each file is expected to be in the VTK XML unstructured grid data format. The standard extension is .vtu. If more than 1 file is specified, the reader will switch to file series mode in which it will pretend that it can support time and provide 1 file per time step. |                  |  |
| <b>Cell Arrays</b><br>( <i>CellArrayStatus</i> )   | This property lists which cell-centered arrays to read.  |                  | The list of array names is provided by the reader. |
| <b>Point Arrays</b><br>( <i>PointArrayStatus</i> ) | This property lists which point-centered arrays to read.   |                  | The list of array names is provided by the reader. |

## XYZ Reader

Read XYZ molecular data files into a polygonal dataset.

The XYZ reader reads XYZ molecular data files. The expected file extension is .xyz. The output of this reader is a polygonal dataset.

| Property                                | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> ) | This property specifies the file name for the XYZ reader.        |                  |              |
| <b>Time Step</b><br>( <i>TimeStep</i> ) | This property specifies the timestep the XYZ reader should load. | 0                |              |

## proSTAR (STARCD) Reader

Reads geometry in proSTAR (STARCD) file format.

ProStarReader creates an unstructured grid dataset.

It reads .cel/.vrt files stored in proSTAR (STARCD) ASCII format.

| Property                                      | Description  | Default Value(s) | Restrictions |
|---|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> )       | Set the file name for the proSTAR (STARCD) reader. |                  |              |
| <b>Scale Factor</b><br>( <i>ScaleFactor</i> ) | Scaling factor for the points                      | 1                |              |

# List of Sources

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## 2D Glyph

Create a 2D glyph (e.g., arrow, cross, dash, etc.)

The 2D Glyph source is used for generating a family of 2D glyphs, each of which lies in the x-y plane. The output of the 2D Glyph source is polygonal data.

| Property                                  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Center</b><br>( <i>Center</i> )        | Set the x, y, z coordinates of the origin of the 2D glyph.  | 0 0 0            |   |
| <b>Filled</b><br>( <i>Filled</i> )        | If the value of this property is 1, the 2D glyph will be a filled polygon; otherwise, only the edges (line segments) will be included in the output. This only applies to closed 2D glyphs. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Glyph Type</b><br>( <i>GlyphType</i> ) | This property specifies the type of the 2D glyph.   | 9                | The value must be one of the following: Vertex (1), Dash (2), Cross (3), ThickCross (4), Triangle (5), Square (6), Circle (7), Diamond (8), Arrow (9), ThickArrow (10), HookedArrow (11), EdgeArrow (12). |

## 3D Text

3D geometric representation of a text string

The 3D Text source displays a text string as polygonal data.

| Property                       | Description   | Default Value(s) | Restrictions |
|--------------------------------|---|------------------|--------------|
| <b>Text</b><br>( <i>Text</i> ) | This property contains the text string to be displayed. The ASCII alphanumeric characters a-z, A-Z, and 0-9 are supported; so are ASCII punctuation marks. The only supported control character is "\n", which inserts a new line in the text string. | 3D Text          |              |

## Annotate Time

Shows the animation time as text annotation in the view.

The Annotate Time source can be used to show the animation time in text annotation.

| Property                           | Description   | Default Value(s) | Restrictions |
|------------------------------------|---|------------------|--------------|
| <b>Format</b><br>( <i>Format</i> ) | This property specifies the format used to display the input time (using printf style). | Time: %f         |              |

---

## Arrow

3D arrow with a long cylindrical shaft and a cone for the tip

The Arrow source appends a cylinder to a cone to form a 3D arrow. The length of the whole arrow is 1.0 unit. The output of the Arrow source is polygonal data. This polygonal data will not contain normals, so rendering of the arrow will be performed using flat shading. The appearance of the arrow can be improved without significantly increasing the resolution of the tip and shaft by generating normals. (Use Normals Generation filter).

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Invert</b><br>( <i>Invert</i> )                    | Inverts the arrow direction.  | 0                | Only the values 0 and 1 are accepted.                                       |
| <b>Shaft Radius</b><br>( <i>ShaftRadius</i> )         | This property specifies the radius of the shaft of the arrow (the cylinder).  | 0.03             | The value must be greater than or equal to 0 and less than or equal to 5.   |
| <b>Shaft Resolution</b><br>( <i>ShaftResolution</i> ) | This property specifies the number of faces of the shaft of the arrow (the cylinder). As the resolution increases, the cylinder will become smoother.             | 6                | The value must be greater than or equal to 0 and less than or equal to 128. |
| <b>Tip Length</b><br>( <i>TipLength</i> )             | This property specifies the length of the tip.  | 0.35             | The value must be greater than or equal to 0 and less than or equal to 1.   |
| <b>Tip Radius</b><br>( <i>TipRadius</i> )             | This property specifies the radius of the widest part of the tip of the arrow (the cone).   | 0.1              | The value must be greater than or equal to 0 and less than or equal to 10.  |
| <b>Tip Resolution</b><br>( <i>TipResolution</i> )     | This property specifies the number of faces in the representation of the tip of the arrow (the cone). As the resolution increases, the cone will become smoother. | 6                | The value must be greater than or equal to 1 and less than or equal to 128. |

## Axes

Three lines representing the axes - red line along X, green line along Y, and blue line along Z

The Axes source can be used to add a representation of the coordinate system axes to the 3D scene. The X axis will be drawn as a blue line, the Y axis as a green line, and the Z axis as a red line. The axes can be drawn either as three lines drawn in the positive direction from the origin or as three lines crossing at the origin (drawn in both the positive and negative directions). The output of the Axes source is polygonal data. This polygonal data has a scalar per line so that the lines can be colored. It also has normals defined.

| Property                                      | Description  | Default Value(s) | Restrictions                                  |
|---|--|------------------|---|
| <b>Origin</b><br>( <i>Origin</i> )            | The values of this property set the X, Y, and Z coordinates of the origin of the axes.   | 0 0 0            |   |
| <b>Scale Factor</b><br>( <i>ScaleFactor</i> ) | By default the axes lines have a length of 1 (or 1 in each direction, for a total length of 2, if value of the Symmetric property is 1). Increasing or decreasing the value of this property will make the axes larger or smaller, respectively. | 1                | The value must be greater than or equal to 0. |
| <b>Symmetric</b><br>( <i>Symmetric</i> )      | When this property is set to 1, the axes extend along each of the positive and negative directions for a distance equal to the value of the Scale Factor property. When set to 0, the axes extend only in the positive direction.                | 0                | Only the values 0 and 1 are accepted.         |

## Box

Create a box with specified X, Y, and Z lengths.

The Box source can be used to add a box to the 3D scene. The output of the Box source is polygonal data containing both normals and texture coordinates.

| Property                     | Description   | Default Value(s) | Restrictions                                  |
|------------------------------|---|------------------|---|
| <b>Center</b><br>(Center)    | This property specifies the center of the box.                    | 0 0 0            |   |
| <b>X Length</b><br>(XLength) | This property specifies the length of the box in the X direction. | 1                | The value must be greater than or equal to 0. |
| <b>Y Length</b><br>(YLength) | This property specifies the length of the box in the Y direction. | 1                | The value must be greater than or equal to 0. |
| <b>Z Length</b><br>(ZLength) | This property specifies the length of the box in the Z direction. | 1                | The value must be greater than or equal to 0. |

## Cone

Create a 3D cone of a given radius and height.

The Cone source can be used to add a polygonal cone to the 3D scene. The output of the Cone source is polygonal data.

| Property                          | Description  | Default Value(s) | Restrictions  |
|-----------------------------------|--|------------------|---|
| <b>Capping</b><br>(Capping)       | If this property is set to 1, the base of the cone will be capped with a filled polygon. Otherwise, the base of the cone will be open.   | 1                | Only the values 0 and 1 are accepted.                                       |
| <b>Center</b><br>(Center)         | This property specifies the center of the cone.  | 0 0 0            |   |
| <b>Direction</b><br>(Direction)   | Set the orientation vector of the cone. The vector does not have to be normalized. The cone will point in the direction specified.   | 1 0 0            |   |
| <b>Height</b><br>(Height)         | This property specifies the height of the cone.  | 1                | The value must be greater than or equal to 0.                               |
| <b>Radius</b><br>(Radius)         | This property specifies the radius of the base of the cone.  | 0.5              | The value must be greater than or equal to 0.                               |
| <b>Resolution</b><br>(Resolution) | This property indicates the number of divisions around the cone. The higher this number, the closer the polygonal approximation will come to representing a cone, and the more polygons it will contain. | 6                | The value must be greater than or equal to 0 and less than or equal to 512. |

## Cylinder

Create a 3D cylinder of a given radius and height.

The Cylinder source can be used to add a polygonal cylinder to the 3D scene. The output of the Cylinder source is polygonal data containing both normals and texture coordinates.

| Property                          | Description  | Default Value(s) | Restrictions  |
|-----------------------------------|--|------------------|---|
| <b>Capping</b><br>(Capping)       | If this property is set to 1, the ends of the cylinder will each be capped with a closed polygon. Otherwise, the ends of the cylinder will be open.  | 1                | Only the values 0 and 1 are accepted.                                       |
| <b>Center</b><br>(Center)         | This property specifies the coordinate value at the center of the cylinder.  | 0 0 0            |   |
| <b>Height</b><br>(Height)         | This property specifies the height of the cylinder (along the y axis).   | 1                | The value must be greater than or equal to 0.                               |
| <b>Radius</b><br>(Radius)         | This property specifies the radius of the cylinder.  | 0.5              | The value must be greater than or equal to 0.                               |
| <b>Resolution</b><br>(Resolution) | This property indicates the number of divisions around the cylinder. The higher this number, the closer the polygonal approximation will come to representing a cylinder, and the more polygons it will contain. | 6                | The value must be greater than or equal to 2 and less than or equal to 512. |

## Data Object Generator

Parses a string to produce composite data objects consisting of simple templated datasets.

vtkDataObjectGenerator parses a string and produces dataobjects from the dataobject template names it sees in the string. For example, if the string contains "ID1" the generator will create a vtkImageData. "UF1", "RG1", "SG1", "PD1", and "UG1" will produce vtkUniformGrid, vtkRectilinearGrid, vtkStructuredGrid, vtkPolyData and vtkUnstructuredGrid respectively. "PD2" will produce an alternate vtkPolydata.

You can compose composite datasets from the atomic ones listed above by placing them within one of the two composite dataset identifiers - "MB{ }" or "HB[ ]". "MB{ ID1 PD1 MB{ } }" for example will create a vtkMultiBlockDataSet consisting of three blocks: image data, poly data, multi-block (empty). Hierarchical Box data sets additionally require the notion of groups, declared within "()" braces, to specify AMR depth. "HB[ (UF1)(UF1)(UF1) ]" will create a vtkHierarchicalBoxDataSet representing an octree that is three levels deep, in which the firstmost cell in each level is refined.

| Property                    | Description  | Default Value(s) | Restrictions |
|-----------------------------|--|------------------|--------------|
| <b>Program</b><br>(Program) | This property contains the string that is parsed to determine the structured of the output data object to produce. | ID1              |              |

## Disk

Create a 3D disk with a specified inner and outer radius.

The Disk source can be used to add a polygonal disk to the 3D scene. The output of the Disk source is polygonal data.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Circumferential Resolution</b><br>( <i>CircumferentialResolution</i> ) | Set the number of points in circumferential direction. | 8                | The value must be greater than or equal to 3 and less than or equal to 1024. |
| <b>Inner Radius</b><br>( <i>InnerRadius</i> )                             | Specify inner radius of hole in disc.                  | 0.5              | The value must be greater than or equal to 0.                                |
| <b>Outer Radius</b><br>( <i>OuterRadius</i> )                             | Specify outer radius of disc.                          | 1                | The value must be greater than or equal to 0.                                |
| <b>Radial Resolution</b><br>( <i>RadialResolution</i> )                   | Set the number of points in radial direction.          | 8                | The value must be greater than or equal to 3 and less than or equal to 1024. |

## Hierarchical Fractal

Test source for AMR with HierarchicalDataSet

The Hierarchical Fractal source is a collection of uniform grids. All have the same dimensions. Each block has a different origin and spacing. This source uses the Mandelbrot source to create cell data. The fractal array is scaled to look like a volume fraction.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Asymetric</b><br>( <i>Asymetric</i> )                | If this property is set to 0, all the blocks will be the same size. If it is set to 1, an asymmetric dataset will be created: some blocks will have an X dimension that is larger by 2 units. | 0                | Only the values 0 and 1 are accepted.                                      |
| <b>Dimensions</b><br>( <i>Dimensions</i> )              | This property specifies the X, Y, Z cell dimensions of a block.   | 10               | The value must be greater than or equal to 2 and less than or equal to 64. |
| <b>Fractal Value</b><br>( <i>FractalValue</i> )         | The value of this property will be mapped to 0.5 for use as a volume fraction.  | 9.5              | The value must be greater than or equal to 0.                              |
| <b>Ghost Levels</b><br>( <i>GhostLevels</i> )           | This property specifies whether ghost levels should be generated at processor boundaries.   | 1                | Only the values 0 and 1 are accepted.                                      |
| <b>Maximum Level</b><br>( <i>MaximumLevel</i> )         | This property specifies how many levels of refinement should be included in this hierarchical dataset.  | 5                | The value must be greater than or equal to 1 and less than or equal to 10. |
| <b>Rectilinear Grids</b><br>( <i>RectilinearGrids</i> ) | If this property is set to 1, the hierarchical dataset will contain rectilinear grids; otherwise it will contain uniform grids.   | 0                | Only the values 0 and 1 are accepted.                                      |
| <b>Time Step</b><br>( <i>TimeStep</i> )                 | This property specifies the timestep to use for this dataset.   | 0                |  |
| <b>Two Dimensional</b><br>( <i>TwoDimensional</i> )     | If this property is set to 1, the generated dataset will be 2D; otherwise it will be 3D.  | 1                | Only the values 0 and 1 are accepted.                                      |

## Line

This source creates a line between two points. The resolution indicates how many segments are in the line.

The Line source can be used to interactively (using a 3D widget) or manually (using the entries on the user interface) add a line to the 3D scene. The output of the Line source is polygonal data.

| Property                                   | Description   | Default Value(s) | Restrictions                                  |
|--|---|------------------|---|
| <b>Point1</b><br>( <i>Point1</i> )         | This property controls the coordinates of the first endpoint of the line.   | -0.5 0 0         |   |
| <b>Point2</b><br>( <i>Point2</i> )         | This property controls the coordinates of the second endpoint of the line.  | 0.5 0 0          |   |
| <b>Resolution</b><br>( <i>Resolution</i> ) | This property specifies the number of pieces into which to divide the line. | 6                | The value must be greater than or equal to 1. |

## Mandelbrot

Representation (unsigned char) of the Mandelbrot set in up to 3 dimensions

The Mandelbrot source can be used to add a uniform rectilinear grid with scalar values derived from the Mandelbrot set to the 3D scene. The equation used is  $z = z^2 + C$  (where  $z$  and  $C$  are complex, and  $C$  is a constant). The scalar values in the grid are the number of iterations of the equation it takes for the magnitude of the value to become greater than 2. In the equation, the initial value of  $z$  is 0. By default, the real component of  $C$  is mapped onto the X axis; the imaginary component of  $C$  is mapped onto the Y axis; and the imaginary component of the initial value is mapped onto the Z axis. If a two-dimensional extent is specified, the resulting image will be displayed. If a three-dimensional extent is used, then the bounding box of the volume will be displayed. The output of the Mandelbrot source is image (uniform rectilinear) data.

| Property  | Description  | Default Value(s)   | Restrictions   |
|---|--|--------------------|--|
| <b>Maximum Number of Iterations</b><br>( <i>MaximumNumberOfIterations</i> ) | The value of this property specifies the limit on computational iterations (i.e., the maximum number of iterations to perform to determine if the value will go above 2). Values less than 2.0 after the specified number of iterations are considered in the fractal set.   | 100                | The value must be greater than or equal to 1 and less than or equal to 5000. |
| <b>Origin CX</b><br>( <i>OriginCX</i> )                                     | The four values of this property indicate (in order) the components of $C$ (real and imaginary) and the components of the initial value, $X$ (real and imaginary).   | -1.75<br>-1.25 0 0 |  |
| <b>Projection Axes</b><br>( <i>ProjectionAxes</i> )                         | The three values in this property allow you to specify the projection from the 4D space used by the Mandelbrot set to the axes of the 3D volume. By default, the real component of $C$ (represented by 0) is mapped to the X axis; the imaginary component of $C$ (represented by 1) is mapped to the Y axis; and the real component of $X$ , the initial value (represented by 2) is mapped to the Z axis. The imaginary component of $X$ is represented by 3. All values entered must be between 0 and 3, inclusive. | 0 1 2              |  |
| <b>Size CX</b><br>( <i>SizeCX</i> )   | The four values of this property indicate the length of the output in each of the four dimensions (the real and imaginary components of $C$ and the real and imaginary components of $X$ ). The three dimensions specified in the Projection Axes property will determine which of these values specify the length of the axes in the output.  | 2.5 2.5 2<br>1.5   |  |
| <b>Subsample Rate</b><br>( <i>SubsampleRate</i> )                           | This property specifies the rate at which to subsample the volume. The extent of the dataset in each dimension will be divided by this value.  | 1                  | The value must be greater than or equal to 1.                                |



|   |  |                    |  |
|---|--|--------------------|--|
| <b>Whole Extent</b><br>( <i>WholeExtent</i> ) | The six values in the property indicate the X, Y, and Z extent of the output data. The first two numbers are the minimum and maximum X extent; the next two are the minimum and maximum Y extent; and the final two are the minimum and maximum Z extent. The numbers are inclusive, so values of 0, 250, 0, 250, 0, 0 indicate that the dimensions of the output will be 251 x 251 x 1. | 0 250 0<br>250 0 0 |  |
|---|--|--------------------|--|

## Octree Fractal

Test source for octree with Mandelbrot fractal

Create an octree from a Mandelbrot fractal. See the Mandelbrot source for a description of the variables used.

| Property  | Description  | Default Value(s)   | Restrictions   |
|---|--|--------------------|--|
| <b>Dimension</b><br>( <i>Dimension</i> )                                    | This property specifies the dimensionality of the fractal: 1D - Binary tree line, 2D - Quadtree plane, 3D - Octree volume.   | 2                  | The value must be greater than or equal to 1 and less than or equal to 3.    |
| <b>Maximum Level</b><br>( <i>MaximumLevel</i> )                             | This property specifies the maximum refinement level for the grid.   | 5                  | The value must be greater than or equal to 2 and less than or equal to 11.   |
| <b>Maximum Number of Iterations</b><br>( <i>MaximumNumberOfIterations</i> ) | The value of this property specifies the limit on computational iterations (i.e., the maximum number of iterations to perform to determine if the value will go above 2). Values less than 2.0 after the specified number of iterations are considered in the fractal set.   | 100                | The value must be greater than or equal to 1 and less than or equal to 5000. |
| <b>Minimum Level</b><br>( <i>MinimumLevel</i> )                             | This property specifies the minimum refinement level for the grid.   | 3                  | The value must be greater than or equal to 1 and less than or equal to 10.   |
| <b>Origin CX</b><br>( <i>OriginCX</i> )                                     | This property specifies the imaginary and real values for C (constant) and X (initial value). See Mandelbrot source for a description of the C and X variables.  | -1.75<br>-1.25 0 0 |  |
| <b>Projection Axes</b><br>( <i>ProjectionAxes</i> )                         | This property indicates which axes of the dataset to display. See Mandelbrot source for a description of the possible axes.  | 0 1 2              |  |
| <b>Size CX</b><br>( <i>SizeCX</i> )   | The four values of this property indicate the length of the output in each of the four dimensions (the real and imaginary components of C and the real and imaginary components of X). The three dimensions specified in the Projection Axes property will determine which of these values specify the length of the axes in the output. | 2.5 2.5 2<br>1.5   |  |
| <b>Threshold</b><br>( <i>Threshold</i> )                                    | This property specifies a threshold value that determines when to subdivide a leaf node.   | 2                  |  |

## Outline

3D outline of the specified bounds.

The Outline source creates an axis aligned bounding box given the user-specified minimum and maximum coordinates for each axis.

| Property                           | Description   | Default Value(s) | Restrictions |
|------------------------------------|---|------------------|--------------|
| <b>Bounds</b><br>( <i>Bounds</i> ) | The values of this property specify the minimum and maximum X, Y, and Z coordinates (X min, X max, Y min, Y max, Z min, Z max) for drawing the outline. | 0 1 0 1 0 1      |              |

## Plane

Create a parallelogram given an origin and two points. The resolution indicates the number of division along each axis of the plane.

The Plane source can be used to add a polygonal parallelogram to the 3D scene. Unlike the sphere, cone, and cylinder sources, the parallelogram is exactly represented at the lowest resolution, but higher resolutions may be desired if this plane is to be used as an input to a filter. The output of the Plane source is polygonal data.

| Property                                      | Description   | Default Value(s) | Restrictions                                  |
|---|---|------------------|---|
| <b>Origin</b><br>( <i>Origin</i> )            | This property specifies the 3D coordinate of the origin (one corner) of the plane.  | -0.5 -0.5 0      |   |
| <b>Point1</b><br>( <i>Point1</i> )            | This property specifies the 3D coordinate a second corner of the parallelogram. The line connecting this point and that specified by the Origin property define one edge of the parallelogram (its X axis).     | 0.5 -0.5 0       |   |
| <b>Point2</b><br>( <i>Point2</i> )            | This property specifies the 3D coordinate a third corner of the parallelogram. The line connecting this point and that specified by the Origin property define a second edge of the parallelogram (its Y axis). | -0.5 0.5 0       |   |
| <b>X Resolution</b><br>( <i>XResolution</i> ) | This property specifies the number of divisions along the X axis of the parallelogram.  | 1                | The value must be greater than or equal to 1. |
| <b>Y Resolution</b><br>( <i>YResolution</i> ) | This property specifies the number of divisions along the Y axis of the parallelogram.  | 1                | The value must be greater than or equal to 1. |

## Point Source

Create a point cloud of a certain size, radius, and center.

The point source creates a specified number of points within a given radius about a specified center point.

| Property   | Description  | Default Value(s) | Restrictions                                  |
|--|--|------------------|---|
| <b>Center</b><br>( <i>Center</i> )                   | This property specifies the 3D coordinates of the center of the point cloud.                           | 0 0 0            |   |
| <b>Number Of Points</b><br>( <i>NumberOfPoints</i> ) | This property specifies the number of points in the point cloud.                                       | 1                | The value must be greater than or equal to 1. |
| <b>Radius</b><br>( <i>Radius</i> )                   | This property specifies the radius of the point cloud, measured from the value of the Center property. | 0                | The value must be greater than or equal to 0. |

## Programmable Source

Executes a user supplied python script to produce an output dataset.

This source will execute a python script to produce an output dataset.

The source keeps a copy of the python script in Script, and creates

Interpreter, a python interpreter to run the script upon the first execution.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Script</b><br>( <b>RequestInformation</b> )<br>( <i>InformationScript</i> ) | This property is a python script that is executed during the RequestInformation pipeline pass. Use this to provide information such as WHOLE_EXTENT to the pipeline downstream. |                  |   |
| <b>Output Data Set Type</b><br>( <i>OutputDataSetType</i> )                    | The value of this property determines the dataset type for the output of the programmable source.   | 0                | The value must be one of the following: Same as Input (8), vtkPolyData (0), vtkStructuredGrid (2), vtkRectilinearGrid (3), vtkUnstructuredGrid (4), vtkImageData (6), vtkMultiblockDataSet (13), vtkHierarchicalBoxDataSet (15), vtkTable (19). |
| <b>Python Path</b><br>( <i>PythonPath</i> )                                    | A semi-colon (;) separated list of directories to add to the python library search path.  |                  |   |
| <b>Script</b><br>( <i>Script</i> )   | This property contains the text of a python program that the programmable source runs.  |                  |   |

## Ruler

This is a line source that can be used to measure distance between two points

The ruler can be used to interactively (using a 3D widget) or manually

(using the entries on the user interface) specify two points and then

determine the distance between the two points. To place points on the

surface of any dataset, one can use the 'p' key shortcut.

| Property                           | Description  | Default Value(s) | Restrictions |
|------------------------------------|--|------------------|--------------|
| <b>Point1</b><br>( <i>Point1</i> ) | This property controls the coordinates of the first endpoint of the line.  | -0.5 0 0         |              |
| <b>Point2</b><br>( <i>Point2</i> ) | This property controls the coordinates of the second endpoint of the line. | 0.5 0 0          |              |

## Sphere

Create a 3D sphere given a center and radius.

The Sphere source can be used to add a polygonal sphere to the 3D scene. The output of the Sphere source is polygonal data with point normals defined.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Center</b><br>( <i>Center</i> )                    | This property specifies the 3D coordinates for the center of the sphere.  | 0 0 0            |   |
| <b>End Phi</b><br>( <i>EndPhi</i> )                   | The value of this property can be adjusted to form only a portion of a sphere. The value is measured in degrees.  | 180              | The value must be greater than or equal to 0 and less than or equal to 180. |
| <b>End Theta</b><br>( <i>EndTheta</i> )               | The value of this property can be adjusted to form only a portion of a sphere. This value is measured in degrees.   | 360              | The value must be greater than or equal to 0 and less than or equal to 360. |
| <b>Phi Resolution</b><br>( <i>PhiResolution</i> )     | The value of this property represents the number of divisions between Start Phi and End Phi on the sphere. (See the Start Phi and End Phi properties.) The phi divisions are similar to latitude lines on the earth.  | 8                | The value must be greater than or equal to 3.                               |
| <b>Radius</b><br>( <i>Radius</i> )                    | This property specifies the radius of the sphere.   | 0.5              | The value must be greater than or equal to 0.                               |
| <b>Start Phi</b><br>( <i>StartPhi</i> )               | To form a complete sphere, the value of this property should be 0 degrees, and the value of the End Phi property should be 180 degrees. The value of this property can be adjusted to form only a portion of a sphere. Set the starting angle (in degrees) in the latitudinal direction.  | 0                | The value must be greater than or equal to 0 and less than or equal to 180. |
| <b>Start Theta</b><br>( <i>StartTheta</i> )           | To form a complete sphere, the value of this property should be 0 degrees, and the value of the End Theta property should be 360 degrees. The value of this property can be adjusted to form only a portion of a sphere.  | 0                | The value must be greater than or equal to 0 and less than or equal to 360. |
| <b>Theta Resolution</b><br>( <i>ThetaResolution</i> ) | The value of this property represents the number of divisions between Start Theta and End Theta around the sphere. (See the Start Theta and End Theta properties.) The theta divisions are similar to longitude lines on the earth. The higher the resolution, the closer the approximation will come to a sphere, and the more polygons there will be. | 8                | The value must be greater than or equal to 3.                               |

## SplineSource

Tessellate parametric functions.

This class tessellates parametric functions. The user must specify how many points in the parametric coordinate directions are required (i.e., the resolution), and the mode to use to generate scalars.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Parametric Function</b><br>( <i>ParametricFunction</i> ) | Property used to reference the parametric function as data generator. |                  | The value must be set to one of the following: Spline. |

## Superquadric

Create a superquadric according to the theta and phi roundness parameters. This one source can generate a wide variety of 3D objects including a box, a sphere, or a torus.

The Superquadric source can be used to add a polygonal superquadric to the 3D scene. This source can be used to create a wide variety of shapes (e.g., a sphere, a box, or a torus) by adjusting the roundness parameters. The output of the Superquadric source is polygonal data with point normals and texture coordinates defined.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Center</b><br>( <i>Center</i> )                    | This property specifies the 3D coordinates of the center of the superquadric.  | 0 0 0            |  |
| <b>Phi Resolution</b><br>( <i>PhiResolution</i> )     | The value of this property represents the number of divisions in the phi (latitudinal) direction. This number will be rounded to the nearest multiple of 4.  | 16               | The value must be greater than or equal to 4 and less than or equal to 1024.   |
| <b>Phi Roundness</b><br>( <i>PhiRoundness</i> )       | This property defines the roundness in the phi (latitudinal) direction. A value of 0 represents a rectangular shape, a value of 1 represents a circular shape, and values greater than 1 produce higher order shapes.  | 1                | The value must be greater than or equal to 0.01.                               |
| <b>Scale</b><br>( <i>Scale</i> )                      | The three values in this property are used to scale the superquadric in X, Y, and Z. The surface normals will be computed correctly even with anisotropic scaling.   | 1 1 1            |  |
| <b>Size</b><br>( <i>Size</i> )                        | The value of this property represents the isotropic size of the superquadric. Note that both the Size and Thickness properties control coefficients of superquadric generation, so the value of this property may not exactly describe the size of the superquadric.   | 0.5              |  |
| <b>Theta Resolution</b><br>( <i>ThetaResolution</i> ) | The value of this property represents the number of divisions in the theta (longitudinal) direction. This value will be rounded to the nearest multiple of 8.  | 16               | The value must be greater than or equal to 8 and less than or equal to 1024.   |
| <b>Theta Roundness</b><br>( <i>ThetaRoundness</i> )   | This property defines the roundness of the superquadric in the theta (longitudinal) direction. A value of 0 represents a rectangular shape, a value of 1 represents a circular shape, and values greater than 1 produce higher order shapes.   | 1                | The value must be greater than or equal to 0.01.                               |
| <b>Thickness</b><br>( <i>Thickness</i> )              | If the value of the Toroidal property is 1, this value represents the thickness of the superquadric as a value between 0 and 1. A value close to 0 leads to a thin object with a large hole, and a value near 1 leads to a thick object with a very small hole. Changing the thickness does not change the outer radius of the superquadric. | 0.3333           | The value must be greater than or equal to 0.0001 and less than or equal to 1. |
| <b>Toroidal</b><br>( <i>Toroidal</i> )                | If the value of this property is 0, the generated superquadric will not contain a hole (i.e., the superquadric will be ellipsoidal). Otherwise, a toroidal object is generated.  | 1                | Only the values 0 and 1 are accepted.  |

## Text

The Text source generates a table containing text.

The Text source is used to generate a 1x1 vtkTable with a single text string.

| Property              | Description                                  | Default Value(s) | Restrictions |
|-----------------------|--|------------------|--------------|
| <b>Text</b><br>(Text) | This property specifies the text to display. | Text             |              |

## Time Source

Produces a single cell uniform grid with data values that vary over a sin(t) wave from t=0 to t=1 (radian).

Produces a single cell uniform grid with data values that vary over a sin(t) wave from t=0 to t=1 (radian).

| Property                            | Description  | Default Value(s) | Restrictions                          |
|-------------------------------------|--|------------------|---------------------------------------|
| <b>Analytic</b><br>(Analytic)       | Makes the time source produce discrete steps of or an analytic sin wave. | 0                | Only the values 0 and 1 are accepted. |
| <b>Growing</b><br>(Growing)         | Makes the time source grow and shrink along Y over time.                 | 0                | Only the values 0 and 1 are accepted. |
| <b>X Amplitude</b><br>(X Amplitude) | Controls how far the data set moves along X over time.                   | 0                |                                       |
| <b>Y Amplitude</b><br>(Y Amplitude) | Controls how far the data set moves along Y over time.                   | 0                |                                       |

## Wavelet

Create a regular rectilinear grid in up to three dimensions with values varying according to a periodic function.

The Wavelet source can be used to create a uniform rectilinear grid in up to three dimensions with values varying according to the following periodic function.

$$OS = M * G * (XM * \sin(XF * x) + YM * \sin(YF * y) + ZM * \cos(ZF * z))$$

OS is the output scalar; M represents the maximum value; G represents the Gaussian; XM, YM, and ZM are the X, Y, and Z magnitude values; and XF, YF, and ZF are the X, Y, and Z frequency values. If a two-dimensional extent is specified, the resulting image will be displayed. If a three-dimensional extent is used, then the bounding box of the volume will be displayed.

| Property   | Description  | Default Value(s) | Restrictions                                  |
|--|--|------------------|---|
| <b>Center</b><br>(Center)                        | This property specifies the 3D coordinates of the center of the dataset.   | 0 0 0            |   |
| <b>Maximum</b><br>(Maximum)                      | This parameter specifies the maximum value (M) of the function.  | 255              |   |
| <b>Standard Deviation</b><br>(StandardDeviation) | This property specifies the standard deviation of the Gaussian used in computing this function.  | 0.5              |   |
| <b>Subsample Rate</b><br>(SubsampleRate)         | This property specifies the rate at which to subsample the volume. The extent of the dataset in each dimension will be divided by this value. (See the Whole Extent property.) | 1                | The value must be greater than or equal to 1. |

|                                      |   |                         |  |
|--------------------------------------|---|-------------------------|--|
| <b>Whole Extent</b><br>(WholeExtent) | The six values in this property indicate the X, Y, and Z extent of the output data. The first two values represent the minimum and maximum X indices, the next two are the minimum and maximum Y indices, and the last two are the minimum and maximum Z indices. | -10 10 -10<br>10 -10 10 |  |
| <b>X Freq</b><br>(XFreq)             | This property specifies the natural frequency in X (XF in the equation).  | 60                      |  |
| <b>X Mag</b><br>(XMag)               | This property specifies the wave amplitude in X (XM in the equation).   | 10                      |  |
| <b>Y Freq</b><br>(YFreq)             | This property specifies the natural frequency in Y (YF in the equation).  | 30                      |  |
| <b>Y Mag</b><br>(YMag)               | This property specifies the wave amplitude in Y (YM in the equation).   | 18                      |  |
| <b>Z Freq</b><br>(ZFreq)             | This property specifies the natural frequency in Z (ZF in the equation).  | 40                      |  |
| <b>Z Mag</b><br>(ZMag)               | This property specifies the wave amplitude in Z (ZM in the equation).   | 5                       |  |

## List of Filters

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### AMR Contour

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Capping</b><br>(Capping)                                      | If this property is on, the the boundary of the data set is capped.   | 1                | Only the values 0 and 1 are accepted.  |
| <b>Isosurface</b><br>(ContourValue)                              | This property specifies the values at which to compute the isosurface.  | 1                | The value must lie within the range of the selected data array.  |
| <b>Degenerate Cells</b><br>(DegenerateCells)                     | If this property is on, a transition mesh between levels is created.  | 1                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>(Input)  | This property specifies the input of the filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkCompositeDataSet. |
| <b>Merge Points</b><br>(MergePoints)                             | Use more memory to merge points on the boundaries of blocks.  | 1                | Only the values 0 and 1 are accepted.  |
| <b>Multiprocess Communication</b><br>(MultiprocessCommunication) | If this property is off, each process executes independantly.   | 1                | Only the values 0 and 1 are accepted.  |
| <b>Contour By</b><br>(SelectInputScalars)                        | This property specifies the name of the cell scalar array from which the contour filter will compute isolines and/or isosurfaces. |                  | An array of scalars is required.   |
| <b>Skip Ghost Copy</b><br>(SkipGhostCopy)                        | A simple test to see if ghost values are already set properly.  | 1                | Only the values 0 and 1 are accepted.  |

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|  |   |   |                                       |
|--|---|---|---------------------------------------|
| <b>Triangulate</b><br>( <i>Triangulate</i> ) | Use triangles instead of quads on capping surfaces. | 1 | Only the values 0 and 1 are accepted. |
|--|---|---|---------------------------------------|

## AMR Dual Clip

Clip with scalars. Tetrahedra.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Degenerate Cells</b><br>( <i>DegenerateCells</i> )                     | If this property is on, a transition mesh between levels is created.                           | 1                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )  | This property specifies the input of the filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkCompositeDataSet. |
| <b>Merge Points</b><br>( <i>MergePoints</i> )                             | Use more memory to merge points on the boundaries of blocks.                                   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Multiprocess Communication</b><br>( <i>MultiprocessCommunication</i> ) | If this property is off, each process executes independantly.                                  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Select Material Arrays</b><br>( <i>SelectMaterialArrays</i> )          | This property specifies the cell arrays from which the clip filter will compute clipped cells. |                  | An array of scalars is required.  |
| <b>Volume Fraction Value</b><br>( <i>VolumeFractionSurfaceValue</i> )     | This property specifies the values at which to compute the isosurface.                         | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.   |

## Annotate Time Filter

Shows input data time as text annnotation in the view.

The Annotate Time filter can be used to show the data time in a text annotation.

| Property                           | Description  | Default Value(s) | Restrictions  |
|------------------------------------|--|------------------|---|
| <b>Format</b><br>( <i>Format</i> ) | The value of this property is a format string used to display the input time. The format string is specified using printf style. | Time: %f         |   |
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the input dataset for which to display the time.   |                  | The selected object must be the result of the following: sources (includes readers), filters. |
| <b>Scale</b><br>( <i>Scale</i> )   | The factor by which the input time is scaled.  | 1                |   |
| <b>Shift</b><br>( <i>Shift</i> )   | The amount of time the input is shifted (after scaling).   | 0                |   |



## Append Attributes

Copies geometry from first input. Puts all of the arrays into the output.

The Append Attributes filter takes multiple input data sets with the same geometry and merges their point and cell attributes to produce a single output containing all the point and cell attributes of the inputs. Any inputs without the same number of points and cells as the first input are ignored. The input data sets must already be collected together, either as a result of a reader that loads multiple parts (e.g., EnSight reader) or because the Group Parts filter has been run to form a collection of data sets.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Append Attributes filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Append Datasets

Takes an input of multiple datasets and output has only one unstructured grid.

The Append Datasets filter operates on multiple data sets of any type (polygonal, structured, etc.). It merges their geometry into a single data set. Only the point and cell attributes that all of the input data sets have in common will appear in the output. The input data sets must already be collected together, either as a result of a reader that loads multiple parts (e.g., EnSight reader) or because the Group Parts filter has been run to form a collection of data sets.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the datasets to be merged into a single dataset by the Append Datasets filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Append Geometry

Takes an input of multiple poly data parts and output has only one part.

The Append Geometry filter operates on multiple polygonal data sets. It merges their geometry into a single data set. Only the point and cell attributes that all of the input data sets have in common will appear in the output.

| Property                         | Description                                  | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | Set the input to the Append Geometry filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |

## Block Scalars

The Level Scalars filter uses colors to show levels of a multiblock dataset.

The Level Scalars filter uses colors to show levels of a multiblock dataset.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Level Scalars filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkMultiBlockDataSet. |

## Calculator

Compute new attribute arrays as function of existing arrays.

The Calculator filter computes a new data array or new point coordinates as a function of existing scalar or vector arrays. If point-centered arrays are used in the computation of a new data array, the resulting array will also be point-centered. Similarly, computations using cell-centered arrays will produce a new cell-centered array. If the function is computing point coordinates, the result of the function must be a three-component vector. The Calculator interface operates similarly to a scientific calculator. In creating the function to evaluate, the standard order of operations applies.

Each of the calculator functions is described below. Unless otherwise noted, enclose the operand in parentheses using the ( and ) buttons.

Clear: Erase the current function (displayed in the read-only text box above the calculator buttons).

/: Divide one scalar by another. The operands for this function are not required to be enclosed in parentheses.

Multiply two scalars, or multiply a vector by a scalar (scalar multiple). The operands for this function are not required to be enclosed in parentheses.

-: Negate a scalar or vector (unary minus), or subtract one scalar or vector from another. The operands for this function are not required to be enclosed in parentheses.

+: Add two scalars or two vectors. The operands for this function are not required to be enclosed in parentheses.

sin: Compute the sine of a scalar.

cos: Compute the cosine of a scalar.

tan: Compute the tangent of a scalar.

asin: Compute the arcsine of a scalar.

acos: Compute the arccosine of a scalar.

atan: Compute the arctangent of a scalar.

sinh: Compute the hyperbolic sine of a scalar.

cosh: Compute the hyperbolic cosine of a scalar.

tanh: Compute the hyperbolic tangent of a scalar.

min: Compute minimum of two scalars.

max: Compute maximum of two scalars.

x^y: Raise one scalar to the power of another scalar. The operands for this function are not required to be enclosed in parentheses.

sqrt: Compute the square root of a scalar.

e^x: Raise e to the power of a scalar.

log: Compute the logarithm of a scalar (deprecated. same as log10).

log10: Compute the logarithm of a scalar to the base 10.

ln: Compute the logarithm of a scalar to the base 'e'.

ceil: Compute the ceiling of a scalar.

floor: Compute the floor of a scalar.

abs: Compute the absolute value of a scalar.

v1.v2: Compute the dot product of two vectors. The operands for this function are not required to be enclosed in parentheses.

cross: Compute cross product of two vectors.

mag: Compute the magnitude of a vector.

norm: Normalize a vector.

The operands are described below.

The digits 0 - 9 and the decimal point are used to enter constant scalar values.

iHat, jHat, and kHat are vector constants representing unit vectors in the X, Y, and Z directions, respectively.

The scalars menu lists the names of the scalar arrays and the components of the vector arrays of either the point-centered or cell-centered data. The vectors menu lists the names of the point-centered or cell-centered vector arrays. The function will be computed for each point (or cell) using the scalar or vector value of the array at that point (or cell).

The filter operates on any type of data set, but the input data set must have at least one scalar or vector array. The arrays can be either point-centered or cell-centered. The Calculator filter's output is of the same data set type as the input.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> )                 | This property determines whether the computation is to be performed on point-centered or cell-centered data.                                    | 0                | The value must be one of the following: point_data (1), cell_data (2), field_data (5).   |
| <b>Coordinate Results</b><br>( <i>CoordinateResults</i> )         | The value of this property determines whether the results of this computation should be used as point coordinates or as a new array.            | 0                | Only the values 0 and 1 are accepted.  |
| <b>Function</b><br>( <i>Function</i> )                            | This property contains the equation for computing the new array.  |                  |  |
| <b>Input</b><br>( <i>Input</i> )                                  | This property specifies the input dataset to the Calculator filter. The scalar and vector variables may be chosen from this dataset's arrays.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Replace Invalid Results</b><br>( <i>ReplaceInvalidValues</i> ) | This property determines whether invalid values in the computation will be replaced with a specific value. (See the ReplacementValue property.) | 1                | Only the values 0 and 1 are accepted.  |
| <b>Replacement Value</b><br>( <i>ReplacementValue</i> )           | If invalid values in the computation are to be replaced with another value, this property contains that value.                                  | 0                |  |
| <b>Result Array Name</b><br>( <i>ResultArrayName</i> )            | This property contains the name for the output array containing the result of this computation.   | Result           |  |

## Cell Centers

Create a point (no geometry) at the center of each input cell.

The Cell Centers filter places a point at the center of each cell in the input data set. The center computed is the parametric center of the cell, not necessarily the geometric or bounding box center. The cell attributes of the input will be associated with these newly created points of the output. You have the option of creating a vertex cell per point in the output. This is useful because vertex cells are rendered, but points are not. The points themselves could be used for placing glyphs (using the Glyph filter). The Cell Centers filter takes any type of data set as input and produces a polygonal data set as output.

| Property                                      | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )              | This property specifies the input to the Cell Centers filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Vertex Cells</b><br>( <i>VertexCells</i> ) | If set to 1, a vertex cell will be generated per point in the output. Otherwise only points will be generated. | 0                | Only the values 0 and 1 are accepted.  |

## Cell Data to Point Data

Create point attributes by averaging cell attributes.

The Cell Data to Point Data filter averages the values of the cell attributes of the cells surrounding a point to compute point attributes. The Cell Data to Point Data filter operates on any type of data set, and the output data set is of the same type as the input.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                    | This property specifies the input to the Cell Data to Point Data filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Pass Cell Data</b><br>( <i>PassCellData</i> )    | If this property is set to 1, then the input cell data is passed through to the output; otherwise, only the generated point data will be available in the output.  | 0                | Only the values 0 and 1 are accepted.  |
| <b>Piece Invariant</b><br>( <i>PieceInvariant</i> ) | If the value of this property is set to 1, this filter will request ghost levels so that the values at boundary points match across processes. NOTE: Enabling this option might cause multiple executions of the data source because more information is needed to remove internal surfaces. | 0                | Only the values 0 and 1 are accepted.  |

## Clean

Merge coincident points if they do not meet a feature edge criteria.

The Clean filter takes polygonal data as input and generates polygonal data as output. This filter can merge duplicate points, remove unused points, and transform degenerate cells into their appropriate forms (e.g., a triangle is converted into a line if two of its points are merged).

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Absolute Tolerance</b><br>( <i>AbsoluteTolerance</i> )         | If merging nearby points (see PointMerging property) and using absolute tolerance (see ToleranceIsAbsolute property), this property specifies the tolerance for performing merging in the spatial units of the input data set.  | 1                | The value must be greater than or equal to 0.   |
| <b>Convert Lines To Points</b><br>( <i>ConvertLinesToPoints</i> ) | If this property is set to 1, degenerate lines (a "line" whose endpoints are at the same spatial location) will be converted to points.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Convert Polys To Lines</b><br>( <i>ConvertPolysToLines</i> )   | If this property is set to 1, degenerate polygons (a "polygon" with only two distinct point coordinates) will be converted to lines.  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Convert Strips To Polys</b><br>( <i>ConvertStripsToPolys</i> ) | If this property is set to 1, degenerate triangle strips (a triangle "strip" containing only one triangle) will be converted to triangles.  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                  | Set the input to the Clean filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Piece Invariant</b><br>( <i>PieceInvariant</i> )               | If this property is set to 1, the whole data set will be processed at once so that cleaning the data set always produces the same results. If it is set to 0, the data set can be processed one piece at a time, so it is not necessary for the entire data set to fit into memory; however the results are not guaranteed to be the same as they would be if the Piece invariant option was on. Setting this option to 0 may produce seams in the output dataset when ParaView is run in parallel. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Point Merging</b><br>( <i>PointMerging</i> )                   | If this property is set to 1, then points will be merged if they are within the specified Tolerance or AbsoluteTolerance (see the Tolerance and AbsoluteTolerance property), depending on the value of the ToleranceIsAbsolute property. (See the ToleranceIsAbsolute property.) If this property is set to 0, points will not be merged.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Tolerance</b><br>( <i>Tolerance</i> )                          | If merging nearby points (see PointMerging property) and not using absolute tolerance (see ToleranceIsAbsolute property), this property specifies the tolerance for performing merging as a fraction of the length of the diagonal of the bounding box of the input data set.   | 0                | The value must be greater than or equal to 0 and less than or equal to 1.   |
| <b>Tolerance Is Absolute</b><br>( <i>ToleranceIsAbsolute</i> )    | This property determines whether to use absolute or relative (a percentage of the bounding box) tolerance when performing point merging.  | 0                | Only the values 0 and 1 are accepted.   |

## Clean to Grid

This filter merges points and converts the data set to unstructured grid.

The Clean to Grid filter merges points that are exactly coincident. It also converts the data set to an unstructured grid. You may wish to do this if you want to apply a filter to your data set that is available for unstructured grids but not for the initial type of your data set (e.g., applying warp vector to volumetric data). The Clean to Grid filter operates on any type of data set.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Clean to Grid filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Clip

Clip with an implicit plane. Clipping does not reduce the dimensionality of the data set. The output data type of this filter is always an unstructured grid.

The Clip filter cuts away a portion of the input data set using an implicit plane. This filter operates on all types of data sets, and it returns unstructured grid data on output.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Clip Type</b><br>( <i>ClipFunction</i> )               | This property specifies the parameters of the clip function (an implicit plane) used to clip the dataset.   |                  | The value must be set to one of the following: Plane, Box, Sphere, Scalar.   |
| <b>Input</b><br>( <i>Input</i> )                          | This property specifies the dataset on which the Clip filter will operate.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Inside Out</b><br>( <i>InsideOut</i> )                 | If this property is set to 0, the clip filter will return that portion of the dataset that lies within the clip function. If set to 1, the portions of the dataset that lie outside the clip function will be returned instead.   | 0                | Only the values 0 and 1 are accepted.  |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )           | If clipping with scalars, this property specifies the name of the scalar array on which to perform the clip operation.  |                  | An array of scalars is required.<br><br>Valid array names will be chosen from point and cell data.   |
| <b>Use Value As Offset</b><br>( <i>UseValueAsOffset</i> ) | If UseValueAsOffset is true, Value is used as an offset parameter to the implicit function. Otherwise, Value is used only when clipping using a scalar array.   | 0                | Only the values 0 and 1 are accepted.  |
| <b>Value</b><br>( <i>Value</i> )                          | If clipping with scalars, this property sets the scalar value about which to clip the dataset based on the scalar array chosen. (See SelectInputScalars.) If clipping with a clip function, this property specifies an offset from the clip function to use in the clipping operation. Neither functionality is currently available in ParaView's user interface. | 0                | The value must lie within the range of the selected data array.  |

## Clip Closed Surface

Clip a polygonal dataset with a plane to produce closed surfaces

This clip filter cuts away a portion of the input polygonal dataset using a plane to generate a new polygonal dataset.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Base Color</b><br>( <i>BaseColor</i> )                       | Specify the color for the faces from the input.   | 0.1 0.1 1        | The value must be greater than or equal to (0, 0, 0) and less than or equal to (1, 1, 1).   |
| <b>Clip Color</b><br>( <i>ClipColor</i> )                       | Specify the color for the capping faces (generated on the clipping interface).  | 1 0.11 0.1       | The value must be greater than or equal to (0, 0, 0) and less than or equal to (1, 1, 1).   |
| <b>Clipping Plane</b><br>( <i>ClippingPlane</i> )               | This property specifies the parameters of the clipping plane used to clip the polygonal data.   |                  | The value must be set to one of the following: Plane.   |
| <b>Generate Cell Origins</b><br>( <i>GenerateColorScalars</i> ) | Generate (cell) data for coloring purposes such that the newly generated cells (including capping faces and clipping outlines) can be distinguished from the input cells.   | 0                | Only the values 0 and 1 are accepted.   |
| <b>Generate Faces</b><br>( <i>GenerateFaces</i> )               | Generate polygonal faces in the output.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Generate Outline</b><br>( <i>GenerateOutline</i> )           | Generate clipping outlines in the output wherever an input face is cut by the clipping plane.   | 0                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                | This property specifies the dataset on which the Clip filter will operate.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Inside Out</b><br>( <i>InsideOut</i> )                       | If this flag is turned off, the clipper will return the portion of the data that lies within the clipping plane. Otherwise, the clipper will return the portion of the data that lies outside the clipping plane. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Clipping Tolerance</b><br>( <i>Tolerance</i> )               | Specify the tolerance for creating new points. A small value might incur degenerate triangles.  | 1e-06            |   |

## Compute Derivatives

This filter computes derivatives of scalars and vectors.

CellDerivatives is a filter that computes derivatives of scalars and vectors at the center of cells. You can choose to generate different output including the scalar gradient (a vector), computed tensor vorticity (a vector), gradient of input vectors (a tensor), and strain matrix of the input vectors (a tensor); or you may choose to pass data through to the output.

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                         | This property specifies the input to the filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Output Tensor Type</b><br>( <i>OutputTensorType</i> ) | This property controls how the filter works to generate tensor cell data. You can choose to compute the gradient of the input vectors, or compute the strain tensor of the vector gradient tensor. By default, the filter will take the gradient of the vector data to construct a tensor. | 1                | The value must be one of the following: Nothing (0), Vector Gradient (1), Strain (2).  |
| <b>Output Vector Type</b><br>( <i>OutputVectorType</i> ) | This property Controls how the filter works to generate vector cell data. You can choose to compute the gradient of the input scalars, or extract the vorticity of the computed vector gradient tensor. By default, the filter will take the gradient of the input scalar data.            | 1                | The value must be one of the following: Nothing (0), Scalar Gradient (1), Vorticity (2).   |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )          | This property indicates the name of the scalar array to differentiate.   |                  | An array of scalars is required.   |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )          | This property indicates the name of the vector array to differentiate.   | 1                | An array of vectors is required.   |

## Connectivity

Mark connected components with integer point attribute array.

The Connectivity filter assigns a region id to connected components of the input data set. (The region id is assigned as a point scalar value.) This filter takes any data set type as input and produces unstructured grid output.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Color Regions</b><br>( <i>ColorRegions</i> )     | Controls the coloring of the connected regions.               | 1                | Only the values 0 and 1 are accepted.   |
| <b>Extraction Mode</b><br>( <i>ExtractionMode</i> ) | Controls the extraction of connected surfaces.                | 5                | The value must be one of the following: Extract Point Seeded Regions (1), Extract Cell Seeded Regions (2), Extract Specified Regions (3), Extract Largest Region (4), Extract All Regions (5), Extract Closes Point Region (6). |
| <b>Input</b><br>( <i>Input</i> )                    | This property specifies the input to the Connectivity filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.                        |

## Contingency Statistics

Compute a statistical model of a dataset and/or assess the dataset with a statistical model.

This filter either computes a statistical model of a dataset or takes such a model as its second input. Then, the model (however it is obtained) may optionally be used to assess the input dataset.

This filter computes contingency tables between pairs of attributes. This result is a tabular bivariate probability distribution which serves as a Bayesian-style prior model. Data is assessed by computing

- the probability of observing both variables simultaneously;
- the probability of each variable conditioned on the other (the two values need not be identical); and



- the pointwise mutual information (PMI).

Finally, the summary statistics include the information entropy of the observations.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> )       | Specify which type of field data the arrays will be drawn from.   | 0                | Valid array names will be chosen from point and cell data.   |
| <b>Input</b><br>( <i>Input</i> )                        | The input to the filter. Arrays from this dataset will be used for computing statistics and/or assessed by a statistical model.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData, vtkStructuredGrid, vtkPolyData, vtkUnstructuredGrid, vtkTable, vtkGraph. |
| <b>Model Input</b><br>( <i>ModelInput</i> )             | A previously-calculated model with which to assess a separate dataset. This input is optional.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable, vtkMultiBlockDataSet.   |
| <b>Variables of Interest</b><br>( <i>SelectArrays</i> ) | Choose arrays whose entries will be used to form observations for statistical analysis.   |                  | An array of scalars is required.   |
| <b>Task</b><br>( <i>Task</i> )                          | Specify the task to be performed: modeling and/or assessment.<br><ol style="list-style-type: none"> <li>1. "Statistics of all the data," creates an output table (or tables) summarizing the <b>entire</b> input dataset;</li> <li>2. "Model a subset of the data," creates an output table (or tables) summarizing a <b>randomly-chosen subset</b> of the input dataset;</li> <li>3. "Assess the data with a model," adds attributes to the first input dataset using a model provided on the second input port; and</li> <li>4. "Model and assess the same data," is really just operations 2 and 3 above applied to the same input dataset. The model is first trained using a fraction of the input data and then the entire dataset is assessed using that model.</li> </ol> <p>When the task includes creating a model (i.e., tasks 2, and 4), you may adjust the fraction of the input dataset used for training. You should avoid using a large fraction of the input data for training as you will then not be able to detect overfitting. The <i>Training fraction</i> setting will be ignored for tasks 1 and 3.</p> | 3                | The value must be one of the following: Statistics of all the data (0), Model a subset of the data (1), Assess the data with a model (2), Model and assess the same data (3).  |
| <b>Training Fraction</b><br>( <i>TrainingFraction</i> ) | Specify the fraction of values from the input dataset to be used for model fitting. The exact set of values is chosen at random from the dataset.   | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Contour

Generate isolines or isosurfaces using point scalars.

The Contour filter computes isolines or isosurfaces using a selected point-centered scalar array. The Contour filter operates on any type of data set, but the input is required to have at least one point-centered scalar (single-component) array. The output of this filter is polygonal.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Compute Gradients</b><br>( <i>ComputeGradients</i> ) | If this property is set to 1, a scalar array containing a gradient value at each point in the isosurface or isoline will be created by this filter; otherwise an array of gradients will not be computed. This operation is fairly expensive both in terms of computation time and memory required, so if the output dataset produced by the contour filter will be processed by filters that modify the dataset's topology or geometry, it may be wise to set the value of this property to 0. Not that if ComputeNormals is set to 1, then gradients will have to be calculated, but they will only be stored in the output dataset if ComputeGradients is also set to 1. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compute Normals</b><br>( <i>ComputeNormals</i> )     | If this property is set to 1, a scalar array containing a normal value at each point in the isosurface or isoline will be created by the contour filter; otherwise an array of normals will not be computed. This operation is fairly expensive both in terms of computation time and memory required, so if the output dataset produced by the contour filter will be processed by filters that modify the dataset's topology or geometry, it may be wise to set the value of this property to 0. Select whether to compute normals.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Compute Scalars</b><br>( <i>ComputeScalars</i> )     | If this property is set to 1, an array of scalars (containing the contour value) will be added to the output dataset. If set to 0, the output will not contain this array.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Isosurfaces</b><br>( <i>ContourValues</i> )          | This property specifies the values at which to compute isosurfaces/isolines and also the number of such values.   |                  | The value must lie within the range of the selected data array.   |
| <b>Input</b><br>( <i>Input</i> )                        | This property specifies the input dataset to be used by the contour filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Point Merge Method</b><br>( <i>Locator</i> )         | This property specifies an incremental point locator for merging duplicate / coincident points.   |                  | The selected object must be the result of the following:<br>incremental_point_locators.<br><br>The value must be set to one of the following: MergePoints, IncrementalOctreeMergePoints, NonMergingPointLocator.  |
| <b>Contour By</b><br>( <i>SelectInputScalars</i> )      | This property specifies the name of the scalar array from which the contour filter will compute isolines and/or isosurfaces.  |                  | An array of scalars is required.<br><br>Valid array names will be chosen from point and cell data.  |

## Curvature

This filter will compute the Gaussian or mean curvature of the mesh at each point.

The Curvature filter computes the curvature at each point in a polygonal data set. This filter supports both Gaussian and mean curvatures.

the type can be selected from the Curvature type menu button.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Curvature Type</b><br>( <i>CurvatureType</i> )              | This property specifies which type of curvature to compute.  | 0                | The value must be one of the following: Gaussian (0), Mean (1).   |
| <b>Input</b><br>( <i>Input</i> )                               | This property specifies the input to the Curvature filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Invert Mean Curvature</b><br>( <i>InvertMeanCurvature</i> ) | If this property is set to 1, the mean curvature calculation will be inverted. This is useful for meshes with inward-pointing normals. | 0                | Only the values 0 and 1 are accepted.   |

## D3

Repartition a data set into load-balanced spatially convex regions. Create ghost cells if requested.

The D3 filter is available when ParaView is run in parallel. It operates on any type of data set to evenly divide it across the processors into spatially contiguous regions. The output of this filter is of type unstructured grid.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Boundary Mode</b><br>( <i>BoundaryMode</i> )      | This property determines how cells that lie on processor boundaries are handled. The "Assign cells uniquely" option assigns each boundary cell to exactly one process, which is useful for isosurfacing. Selecting "Duplicate cells" causes the cells on the boundaries to be copied to each process that shares that boundary. The "Divide cells" option breaks cells across process boundary lines so that pieces of the cell lie in different processes. This option is useful for volume rendering. | 0                | The value must be one of the following: Assign cells uniquely (0), Duplicate cells (1), Divide cells (2).  |
| <b>Input</b><br>( <i>Input</i> )                     | This property specifies the input to the D3 filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Minimal Memory</b><br>( <i>UseMinimalMemory</i> ) | If this property is set to 1, the D3 filter requires communication routines to use minimal memory than without this restriction.  | 0                | Only the values 0 and 1 are accepted.  |

## Decimate

Simplify a polygonal model using an adaptive edge collapse algorithm. This filter works with triangles only.

The Decimate filter reduces the number of triangles in a polygonal data set. Because this filter only operates on triangles, first run the Triangulate filter on a dataset that contains polygons other than triangles.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Boundary Vertex Deletion</b><br>( <i>BoundaryVertexDeletion</i> ) | If this property is set to 1, then vertices on the boundary of the dataset can be removed. Setting the value of this property to 0 preserves the boundary of the dataset, but it may cause the filter not to reach its reduction target.                                    | 1                | Only the values 0 and 1 are accepted.   |
| <b>Feature Angle</b><br>( <i>FeatureAngle</i> )                      | The value of this property is used in determining where the data set may be split. If the angle between two adjacent triangles is greater than or equal to the FeatureAngle value, then their boundary is considered a feature edge where the dataset can be split.         | 15               | The value must be greater than or equal to 0 and less than or equal to 180.   |
| <b>Input</b><br>( <i>Input</i> )                                     | This property specifies the input to the Decimate filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Preserve Topology</b><br>( <i>PreserveTopology</i> )              | If this property is set to 1, decimation will not split the dataset or produce holes, but it may keep the filter from reaching the reduction target. If it is set to 0, better reduction can occur (reaching the reduction target), but holes in the model may be produced. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Target Reduction</b><br>( <i>TargetReduction</i> )                | This property specifies the desired reduction in the total number of polygons in the output dataset. For example, if the TargetReduction value is 0.9, the Decimate filter will attempt to produce an output dataset that is 10% the size of the input.)                    | 0.9              | The value must be greater than or equal to 0 and less than or equal to 1.   |

## Delaunay 2D

Create 2D Delaunay triangulation of input points. It expects a vtkPointSet as input and produces vtkPolyData as output. The points are expected to be in a mostly planar distribution.

Delaunay2D is a filter that constructs a 2D Delaunay triangulation from a list of input points. These points may be represented by any dataset of type vtkPointSet and subclasses. The output of the filter is a polygonal dataset containing a triangle mesh.

The 2D Delaunay triangulation is defined as the triangulation that satisfies the Delaunay criterion for n-dimensional simplexes (in this case n=2 and the simplexes are triangles). This criterion states that a circumsphere of each simplex in a triangulation contains only the n+1 defining points of the simplex. In two dimensions, this translates into an optimal triangulation. That is, the maximum interior angle of any triangle is less than or equal to that of any possible triangulation.

Delaunay triangulations are used to build topological structures from unorganized (or unstructured) points. The input to this filter is a list of points specified in 3D, even though the triangulation is 2D. Thus the triangulation is constructed in the x-y plane, and the z coordinate is ignored (although carried through to the output). You can use the option ProjectionPlaneMode in order to compute the best-fitting plane to the set of points, project the points and that plane and then perform the triangulation using their projected positions and then use it as the plane in which the triangulation is performed.

The Delaunay triangulation can be numerically sensitive in some cases. To prevent problems, try to avoid injecting points that will result in triangles with bad aspect ratios (1000:1 or greater). In practice this means inserting points that are "widely dispersed", and enables smooth transition of triangle sizes throughout the mesh. (You may even want to add extra points to create a better point distribution.) If numerical problems are present, you will see a warning message to this effect at the end of the triangulation process.

Warning:

Points arranged on a regular lattice (termed degenerate cases) can be triangulated in more than one way (at least according to the Delaunay criterion). The choice of triangulation (as implemented by this algorithm) depends on the order of the input points. The first three points will form a triangle; other degenerate points will not break this triangle.

Points that are coincident (or nearly so) may be discarded by the algorithm. This is because the Delaunay triangulation requires unique input points. The output of the Delaunay triangulation is supposedly a convex hull. In certain cases this implementation may not generate the convex hull.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Alpha</b><br>( <i>Alpha</i> )                                  | The value of this property controls the output of this filter. For a non-zero alpha value, only edges or triangles contained within a sphere centered at mesh vertices will be output. Otherwise, only triangles will be output.                   | 0                | The value must be greater than or equal to 0.   |
| <b>Bounding Triangulation</b><br>( <i>BoundingTriangulation</i> ) | If this property is set to 1, bounding triangulation points (and associated triangles) are included in the output. These are introduced as an initial triangulation to begin the triangulation process. This feature is nice for debugging output. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                  | This property specifies the input dataset to the Delaunay 2D filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |
| <b>Offset</b><br>( <i>Offset</i> )                                | This property is a multiplier to control the size of the initial, bounding Delaunay triangulation.   | 1                | The value must be greater than or equal to 0.75.  |
| <b>Projection Plane Mode</b><br>( <i>ProjectionPlaneMode</i> )    | This property determines type of projection plane to use in performing the triangulation.  | 0                | The value must be one of the following: XY Plane (0), Best-Fitting Plane (2).   |
| <b>Tolerance</b><br>( <i>Tolerance</i> )                          | This property specifies a tolerance to control discarding of closely spaced points. This tolerance is specified as a fraction of the diagonal length of the bounding box of the points.  | 1e-05            | The value must be greater than or equal to 0 and less than or equal to 1.   |

## Delaunay 3D

Create a 3D Delaunay triangulation of input points. It expects a vtkPointSet as input and produces vtkUnstructuredGrid as output.

Delaunay3D is a filter that constructs a 3D Delaunay triangulation from a list of input points. These points may be represented by any dataset of type vtkPointSet and subclasses. The output of the filter is an unstructured grid dataset. Usually the output is a tetrahedral mesh, but if a non-zero alpha distance value is specified (called the "alpha" value), then only tetrahedra, triangles, edges, and

vertices lying within the alpha radius are output. In other words, non-zero alpha values may result in arbitrary combinations of tetrahedra, triangles, lines, and vertices. (The notion of alpha value is derived from Edelsbrunner's work on "alpha shapes".) The 3D Delaunay triangulation is defined as the triangulation that satisfies the Delaunay criterion for n-dimensional simplexes (in this case  $n=3$  and the simplexes are tetrahedra). This criterion states that a circumsphere of each simplex in a triangulation contains only the  $n+1$  defining points of the simplex. (See text for more information.) While in two dimensions this translates into an "optimal" triangulation, this is not true in 3D, since a measurement for optimality in 3D is not agreed on.

Delaunay triangulations are used to build topological structures from unorganized (or unstructured) points. The input to this filter is a list of points specified in 3D. (If you wish to create 2D triangulations see Delaunay2D.) The output is an unstructured grid.

The Delaunay triangulation can be numerically sensitive. To prevent problems, try to avoid injecting points that will result in triangles with bad aspect ratios (1000:1 or greater). In practice this means inserting points that are "widely dispersed", and enables smooth transition of triangle sizes throughout the mesh. (You may even want to add extra points to create a better point distribution.) If numerical problems are present, you will see a warning message to this effect at the end of the triangulation process.

Warning:

Points arranged on a regular lattice (termed degenerate cases) can be triangulated in more than one way (at least according to the Delaunay criterion). The choice of triangulation (as implemented by this algorithm) depends on the order of the input points. The first four points will form a tetrahedron; other degenerate points (relative to this initial tetrahedron) will not break it.

Points that are coincident (or nearly so) may be discarded by the algorithm. This is because the Delaunay triangulation requires unique input points. You can control the definition of coincidence with the "Tolerance" instance variable.

The output of the Delaunay triangulation is supposedly a convex hull. In certain cases this implementation may not generate the convex hull. This behavior can be controlled by the Offset instance variable. Offset is a multiplier used to control the size of the initial triangulation. The larger the offset value, the more likely you will generate a convex hull; and the more likely you are to see numerical problems.

The implementation of this algorithm varies from the 2D Delaunay algorithm (i.e., Delaunay2D) in an important way. When points are injected into the triangulation, the search for the enclosing

tetrahedron is quite different. In the 3D case, the closest previously inserted point is found, and then the connected tetrahedra are searched to find the containing one. (In 2D, a "walk" towards the enclosing triangle is performed.) If the triangulation is Delaunay, then an enclosing tetrahedron will be found. However, in degenerate cases an enclosing tetrahedron may not be found and the point will be rejected.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Alpha</b><br>( <i>Alpha</i> )                                  | This property specifies the alpha (or distance) value to control the output of this filter. For a non-zero alpha value, only edges, faces, or tetra contained within the circumsphere (of radius alpha) will be output. Otherwise, only tetrahedra will be output. | 0                | The value must be greater than or equal to 0.   |
| <b>Bounding Triangulation</b><br>( <i>BoundingTriangulation</i> ) | This boolean controls whether bounding triangulation points (and associated triangles) are included in the output. (These are introduced as an initial triangulation to begin the triangulation process. This feature is nice for debugging output.)               | 0                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                  | This property specifies the input dataset to the Delaunay 3D filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |
| <b>Offset</b><br>( <i>Offset</i> )                                | This property specifies a multiplier to control the size of the initial, bounding Delaunay triangulation.  | 2.5              | The value must be greater than or equal to 2.5.   |
| <b>Tolerance</b><br>( <i>Tolerance</i> )                          | This property specifies a tolerance to control discarding of closely spaced points. This tolerance is specified as a fraction of the diagonal length of the bounding box of the points.  | 0.001            | The value must be greater than or equal to 0 and less than or equal to 1.   |

## Descriptive Statistics

Compute a statistical model of a dataset and/or assess the dataset with a statistical model.

This filter either computes a statistical model of a dataset or takes such a model as its second input. Then, the model (however it is obtained) may optionally be used to assess the input dataset.

This filter computes the min, max, mean, raw moments M2 through M4, standard deviation, skewness, and kurtosis for each array you select.

The model is simply a univariate Gaussian distribution with the mean and standard deviation provided. Data is assessed using this model by detrending the data (i.e., subtracting the mean) and then dividing by the standard deviation. Thus the assessment is an array whose entries are the number of standard deviations from the mean that each input point lies.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> )          | Specify which type of field data the arrays will be drawn from.   | 0                | Valid array names will be chosen from point and cell data.   |
| <b>Input</b><br>( <i>Input</i> )                           | The input to the filter. Arrays from this dataset will be used for computing statistics and/or assessed by a statistical model.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData, vtkStructuredGrid, vtkPolyData, vtkUnstructuredGrid, vtkTable, vtkGraph. |
| <b>Model Input</b><br>( <i>ModelInput</i> )                | A previously-calculated model with which to assess a separate dataset. This input is optional.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable, vtkMultiBlockDataSet.   |
| <b>Variables of Interest</b><br>( <i>SelectArrays</i> )    | Choose arrays whose entries will be used to form observations for statistical analysis.   |                  | An array of scalars is required.   |
| <b>Deviations should be</b><br>( <i>SignedDeviations</i> ) | Should the assessed values be signed deviations or unsigned?  | 0                | The value must be one of the following: Unsigned (0), Signed (1).  |
| <b>Task</b><br>( <i>Task</i> )                             | Specify the task to be performed: modeling and/or assessment.<br><ol style="list-style-type: none"> <li>"Statistics of all the data," creates an output table (or tables) summarizing the <b>entire</b> input dataset;</li> <li>"Model a subset of the data," creates an output table (or tables) summarizing a <b>randomly-chosen subset</b> of the input dataset;</li> <li>"Assess the data with a model," adds attributes to the first input dataset using a model provided on the second input port; and</li> <li>"Model and assess the same data," is really just operations 2 and 3 above applied to the same input dataset. The model is first trained using a fraction of the input data and then the entire dataset is assessed using that model.</li> </ol> <p>When the task includes creating a model (i.e., tasks 2, and 4), you may adjust the fraction of the input dataset used for training. You should avoid using a large fraction of the input data for training as you will then not be able to detect overfitting. The <i>Training fraction</i> setting will be ignored for tasks 1 and 3.</p> | 3                | The value must be one of the following: Statistics of all the data (0), Model a subset of the data (1), Assess the data with a model (2), Model and assess the same data (3).  |
| <b>Training Fraction</b><br>( <i>TrainingFraction</i> )    | Specify the fraction of values from the input dataset to be used for model fitting. The exact set of values is chosen at random from the dataset.   | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.  |



## Elevation

Create point attribute array by projecting points onto an elevation vector.

The Elevation filter generates point scalar values for an input dataset along a specified direction vector.

The Input menu allows the user to select the data set to which this filter will be applied. Use the Scalar range entry boxes to specify the minimum and maximum scalar value to be generated. The Low Point and High Point define a line onto which each point of the data set is projected. The minimum scalar value is associated with the Low Point, and the maximum scalar value is associated with the High Point. The scalar value for each point in the data set is determined by the location along the line to which that point projects.

| Property                                      | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>High Point</b><br>( <i>HighPoint</i> )     | This property defines the other end of the direction vector (large scalar values). | 0 0 1            | The coordinate must lie within the bounding box of the dataset. It will default to the maximum in each dimension.  |
| <b>Input</b><br>( <i>Input</i> )              | This property specifies the input dataset to the Elevation filter.                 |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Low Point</b><br>( <i>LowPoint</i> )       | This property defines one end of the direction vector (small scalar values).       | 0 0 0            | The coordinate must lie within the bounding box of the dataset. It will default to the minimum in each dimension.  |
| <b>Scalar Range</b><br>( <i>ScalarRange</i> ) | This property determines the range into which scalars will be mapped.              | 0 1              |  |

## Extract AMR Blocks

This filter extracts a list of datasets from hierarchical datasets.

This filter extracts a list of datasets from hierarchical datasets.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                         | This property specifies the input to the Extract Datasets filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkHierarchicalBoxDataSet. |
| <b>Selected Data Sets</b><br>( <i>SelectedDataSets</i> ) | This property provides a list of datasets to extract.             |                  |   |

## Extract Block

This filter extracts a range of blocks from a multiblock dataset.

This filter extracts a range of groups from a multiblock dataset

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Block Indices</b><br>( <i>BlockIndices</i> )           | This property lists the ids of the blocks to extract from the input multiblock dataset.  |                  |  |
| <b>Input</b><br>( <i>Input</i> )                          | This property specifies the input to the Extract Group filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkMultiBlockDataSet. |
| <b>Maintain Structure</b><br>( <i>MaintainStructure</i> ) | This is used only when PruneOutput is ON. By default, when pruning the output i.e. remove empty blocks, if node has only 1 non-null child block, then that node is removed. To preserve these parent nodes, set this flag to true. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Prune Output</b><br>( <i>PruneOutput</i> )             | When set, the output multiblock dataset will be pruned to remove empty nodes. On by default.   | 1                | Only the values 0 and 1 are accepted.  |

## Extract CTH Parts

Create a surface from a CTH volume fraction.

Extract CTH Parts is a specialized filter for visualizing the data from a CTH simulation. It first converts the selected cell-centered arrays to point-centered ones. It then contours each array at a value of 0.5. The user has the option of clipping the resulting surface(s) with a plane. This filter only operates on unstructured data. It produces polygonal output.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Double Volume Arrays</b><br>( <i>AddDoubleVolumeArrayName</i> )                   | This property specifies the name(s) of the volume fraction array(s) for generating parts.   |                  | An array of scalars is required.   |
| <b>Float Volume Arrays</b><br>( <i>AddFloatVolumeArrayName</i> )                     | This property specifies the name(s) of the volume fraction array(s) for generating parts.   |                  | An array of scalars is required.   |
| <b>Unsigned Character Volume Arrays</b><br>( <i>AddUnsignedCharVolumeArrayName</i> ) | This property specifies the name(s) of the volume fraction array(s) for generating parts.   |                  | An array of scalars is required.   |
| <b>Clip Type</b><br>( <i>ClipPlane</i> )   | This property specifies whether to clip the dataset, and if so, it also specifies the parameters of the plane with which to clip. |                  | The value must be set to one of the following: None, Plane, Box, Sphere.   |
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the input to the Extract CTH Parts filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a cell array with 1 components.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Volume Fraction Value</b><br>( <i>VolumeFractionSurfaceValue</i> )                | The value of this property is the volume fraction value for the surface.  | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Extract Cells By Region

This filter extracts cells that are inside/outside a region or at a region boundary.

This filter extracts from its input dataset all cells that are either completely inside or outside of a specified region (implicit function). On output, the filter generates an unstructured grid.

To use this filter you must specify a region (implicit function). You must also specify whether to extract cells lying inside or outside of the region. An option exists to extract cells that are neither inside or outside (i.e., boundary).

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Extract intersected</b><br>( <i>Extract intersected</i> )           | This parameter controls whether to extract cells that are on the boundary of the region.   | 0                | Only the values 0 and 1 are accepted.  |
| <b>Extract only intersected</b><br>( <i>Extract only intersected</i> ) | This parameter controls whether to extract only cells that are on the boundary of the region. If this parameter is set, the Extraction Side parameter is ignored. If Extract Intersected is off, this parameter has no effect. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Extraction Side</b><br>( <i>ExtractInside</i> )                     | This parameter controls whether to extract cells that are inside or outside the region.  | 1                | The value must be one of the following: outside (0), inside (1).   |
| <b>Intersect With</b><br>( <i>ImplicitFunction</i> )                   | This property sets the region used to extract cells.   |                  | The value must be set to one of the following: Plane, Box, Sphere.   |
| <b>Input</b><br>( <i>Input</i> )                                       | This property specifies the input to the Slice filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Extract Edges

Extract edges of 2D and 3D cells as lines.

The Extract Edges filter produces a wireframe version of the input dataset by extracting all the edges of the dataset's cells as lines. This filter operates on any type of data set and produces polygonal output.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Extract Edges filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Extract Level

This filter extracts a range of groups from a hierarchical dataset.

This filter extracts a range of levels from a hierarchical dataset

| Property                           | Description  | Default Value(s) | Restrictions  |
|------------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the input to the Extract Group filter.                 |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkHierarchicalBoxDataSet. |
| <b>Levels</b><br>( <i>Levels</i> ) | This property lists the levels to extract from the input hierarchical dataset. |                  |   |

## Extract Selection

Extract different type of selections.

This filter extracts a set of cells/points given a selection.

The selection can be obtained from a rubber-band selection (either cell, visible or in a frustum) or threshold selection and passed to the filter or specified by providing an ID list.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                        | This property specifies the input from which the selection is extracted.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet, vtkTable. |
| <b>Preserve Topology</b><br>( <i>PreserveTopology</i> ) | If this property is set to 1 the output preserves the topology of its input and adds an insidedness array to mark which cells are inside or out. If 0 then the output is an unstructured grid which contains only the subset of cells that are inside. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Selection</b><br>( <i>Selection</i> )                | The input that provides the selection object.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkSelection.         |
| <b>Show Bounds</b><br>( <i>ShowBounds</i> )             | For frustum selection, if this property is set to 1 the output is the outline of the frustum instead of the contents of the input that lie within the frustum.   | 0                | Only the values 0 and 1 are accepted.  |

## Extract Subset

Extract a subgrid from a structured grid with the option of setting subsample strides.

The Extract Grid filter returns a subgrid of a structured input data set (uniform rectilinear, curvilinear, or nonuniform rectilinear). The output data set type of this filter is the same as the input type.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Include Boundary</b><br>( <i>IncludeBoundary</i> ) | If the value of this property is 1, then if the sample rate in any dimension is greater than 1, the boundary indices of the input dataset will be passed to the output even if the boundary extent is not an even multiple of the sample rate in a given dimension. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )                      | This property specifies the input to the Extract Grid filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkImageData, vtkRectilinearGrid,<br>vtkStructuredPoints, vtkStructuredGrid. |
| <b>Sample Rate I</b><br>( <i>SampleRateI</i> )        | This property indicates the sampling rate in the I dimension. A value greater than 1 results in subsampling; every nth index will be included in the output.  | 1                | The value must be greater than or equal to 1.  |
| <b>Sample Rate J</b><br>( <i>SampleRateJ</i> )        | This property indicates the sampling rate in the J dimension. A value greater than 1 results in subsampling; every nth index will be included in the output.  | 1                | The value must be greater than or equal to 1.  |
| <b>Sample Rate K</b><br>( <i>SampleRateK</i> )        | This property indicates the sampling rate in the K dimension. A value greater than 1 results in subsampling; every nth index will be included in the output.  | 1                | The value must be greater than or equal to 1.  |
| <b>V OI</b><br>( <i>VOI</i> )                         | This property specifies the minimum and maximum point indices along each of the I, J, and K axes; these values indicate the volume of interest (VOI). The output will have the (I,J,K) extent specified here.   | 0 0 0 0 0        | The values must lie within the extent of the input dataset.  |

## Extract Surface

Extract a 2D boundary surface using neighbor relations to eliminate internal faces.

The Extract Surface filter extracts the polygons forming the outer surface of the input dataset. This filter operates on any type of data and produces polygonal data as output.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Extract Surface filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

|  |   |   |   |
|--|---|---|---|
| <b>Nonlinear Subdivision Level</b><br>( <i>NonlinearSubdivisionLevel</i> ) | If the input is an unstructured grid with nonlinear faces, this parameter determines how many times the face is subdivided into linear faces. If 0, the output is the equivalent of its linear counterpart (and the midpoints determining the nonlinear interpolation are discarded). If 1, the nonlinear face is triangulated based on the midpoints. If greater than 1, the triangulated pieces are recursively subdivided to reach the desired subdivision. Setting the value to greater than 1 may cause some point data to not be passed even if no quadratic faces exist. This option has no effect if the input is not an unstructured grid. | 1 | The value must be greater than or equal to 0 and less than or equal to 4. |
| <b>Piece Invariant</b><br>( <i>PieceInvariant</i> )                        | If the value of this property is set to 1, internal surfaces along process boundaries will be removed. NOTE: Enabling this option might cause multiple executions of the data source because more information is needed to remove internal surfaces.  | 1 | Only the values 0 and 1 are accepted.                                     |

## FFT Of Selection Over Time

Extracts selection over time and plots the FFT

Extracts the data of a selection (e.g. points or cells) over time, takes the FFT of them, and plots them.

| Property                                 | Description                                      | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )         | The input from which the selection is extracted. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet, vtkTable, vtkCompositeDataSet. |
| <b>Selection</b><br>( <i>Selection</i> ) | The input that provides the selection object.    |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkSelection.                              |

## FOF/SOD Halo Finder

Sorry, no help is currently available.

| Property   | Description   | Default Value(s) | Restrictions                                  |
|--|---|------------------|---|
| <b>bb (linking length)</b><br>( <i>BB</i> )  | Linking length measured in units of interparticle spacing and is dimensionless. Used to link particles into halos for the friends-of-friends (FOF) algorithm. | 0.2              | The value must be greater than or equal to 0. |
| <b>Compute the most bound particle</b><br>( <i>ComputeMostBoundParticle</i> )            | If checked, the most bound particle for an FOF halo will be calculated. WARNING: This can be very slow.   | 0                | Only the values 0 and 1 are accepted.         |
| <b>Compute the most connected particle</b><br>( <i>ComputeMostConnectedParticle</i> )    | If checked, the most connected particle for an FOF halo will be calculated. WARNING: This can be very slow.   | 0                | Only the values 0 and 1 are accepted.         |
| <b>Compute spherical overdensity (SOD) halos</b><br>( <i>ComputeSOD</i> )                | If checked, spherical overdensity (SOD) halos will be calculated in addition to friends-of-friends (FOF) halos.   | 0                | Only the values 0 and 1 are accepted.         |
| <b>Copy FOF halo catalog to original particles</b><br>( <i>CopyHaloDataToParticles</i> ) | If checked, the friends-of-friends (FOF) halo catalog information will be copied to the original particles as well.   | 0                | Only the values 0 and 1 are accepted.         |

|  |  |     |   |
|--|--|-----|---|
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the input of the filter.   |     | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |
| <b>np (number of seeded particles in one dimension, i.e., total particles = np^3)</b><br>( <i>NP</i> ) | Number of seeded particles in one dimension. Therefore, total simulation particles is np^3 (cubed).  | 256 | The value must be greater than or equal to 0.   |
| <b>overlap (shared point/ghost cell gap distance)</b><br>( <i>Overlap</i> )                            | The space (in rL units) to extend processor particle ownership for ghost particles/cells. Needed for correct halo calculation when halos cross processor boundaries in parallel computation. | 5   | The value must be greater than or equal to 0.   |
| <b>pmin (minimum particle threshold for an FOF halo)</b><br>( <i>PMin</i> )                            | Minimum number of particles (threshold) needed before a group is called a friends-of-friends (FOF) halo.   | 100 | The value must be greater than or equal to 1.   |
| <b>rL (physical box side length)</b><br>( <i>RL</i> )  | The box side length used to wrap particles around if they exceed rL (or less than 0) in any dimension (only positive positions are allowed in the input, or they are wrapped around).        | 100 | The value must be greater than or equal to 0.   |
| <b>scale factor for rho_c</b><br>( <i>RhoCScale</i> )  | Scale factor for rho_c in SOD halo finding such that rho_c' = rho_c * scale factor. Initial rho_c is 2.77536627e11 (M_sun/h) / (Mpc/h)^3.  | 1   |   |
| <b>initial SOD center</b><br>( <i>SODCenterType</i> )  | The initial friends-of-friends (FOF) center used for calculating a spherical overdensity (SOD) halo. WARNING: Using MBP or MCP can be very slow.   | 0   | The value must be one of the following: Center of mass (0), Average position (1), Most bound particle (2), Most connected particle (3).   |
| <b>scale factor for initial SOD mass</b><br>( <i>SODMassScale</i> )                                    | Scale factor for the initial SOD mass such that mass' = mass * scale factor. Initial SOD mass is 1.0e14 (M_sun/h).   | 1   |   |

## Feature Edges

This filter will extract edges along sharp edges of surfaces or boundaries of surfaces.

The Feature Edges filter extracts various subsets of edges from the input data set. This filter operates on polygonal data and produces polygonal output.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Boundary Edges</b><br>( <i>BoundaryEdges</i> ) | If the value of this property is set to 1, boundary edges will be extracted. Boundary edges are defined as lines cells or edges that are used by only one polygon.   | 1                | Only the values 0 and 1 are accepted.                                       |
| <b>Coloring</b><br>( <i>Coloring</i> )            | If the value of this property is set to 1, then the extracted edges are assigned a scalar value based on the type of the edge.   | 0                | Only the values 0 and 1 are accepted.                                       |
| <b>Feature Angle</b><br>( <i>FeatureAngle</i> )   | This value of this property is used to define a feature edge. If the surface normal between two adjacent triangles is at least as large as this Feature Angle, a feature edge exists. (See the FeatureEdges property.) | 30               | The value must be greater than or equal to 0 and less than or equal to 180. |

|  |  |   |   |
|--|--|---|---|
| <b>Feature Edges</b><br>( <i>FeatureEdges</i> )          | If the value of this property is set to 1, feature edges will be extracted. Feature edges are defined as edges that are used by two polygons whose dihedral angle is greater than the feature angle. (See the FeatureAngle property.) Toggle whether to extract feature edges. | 1 | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                         | This property specifies the input to the Feature Edges filter.   |   | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Manifold Edges</b><br>( <i>ManifoldEdges</i> )        | If the value of this property is set to 1, manifold edges will be extracted. Manifold edges are defined as edges that are used by exactly two polygons.  | 0 | Only the values 0 and 1 are accepted.   |
| <b>Non-Manifold Edges</b><br>( <i>NonManifoldEdges</i> ) | If the value of this property is set to 1, non-manifold edges will be extracted. Non-manifold edges are defined as edges that are use by three or more polygons.   | 1 | Only the values 0 and 1 are accepted.   |

## Generate Ids

Generate scalars from point and cell ids.

This filter generates scalars using cell and point ids. That is, the point attribute data scalars are generated from the point ids, and the cell attribute data scalars or field data are generated from the the cell ids.

| Property                                  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Array Name</b><br>( <i>ArrayName</i> ) | The name of the array that will contain ids.                             | Ids              |  |
| <b>Input</b><br>( <i>Input</i> )          | This property specifies the input to the Cell Data to Point Data filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Generate Quadrature Points

Create a point set with data at quadrature points.

"Create a point set with data at quadrature points."

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                           | This property specifies the input of the filter.                     |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |
| <b>Select Source Array</b><br>( <i>SelectSourceArray</i> ) | Specifies the offset array from which we generate quadrature points. |                  | An array of scalars is required.  |



## Generate Quadrature Scheme Dictionary

Generate quadrature scheme dictionaries in data sets that do not have them.

Generate quadrature scheme dictionaries in data sets that do not have them.

| Property                         | Description                                      | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input of the filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |

## Generate Surface Normals

This filter will produce surface normals used for smooth shading. Splitting is used to avoid smoothing across feature edges.

This filter generates surface normals at the points of the input polygonal dataset to provide smooth shading of the dataset. The resulting dataset is also polygonal. The filter works by calculating a normal vector for each polygon in the dataset and then averaging the normals at the shared points.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Compute Cell Normals</b><br>( <i>ComputeCellNormals</i> )     | This filter computes the normals at the points in the data set. In the process of doing this it computes polygon normals too. If you want these normals to be passed to the output of this filter, set the value of this property to 1.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Consistency</b><br>( <i>Consistency</i> )                     | The value of this property controls whether consistent polygon ordering is enforced. Generally the normals for a data set should either all point inward or all point outward. If the value of this property is 1, then this filter will reorder the points of cells that whose normal vectors are oriented the opposite direction from the rest of those in the data set. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Feature Angle</b><br>( <i>FeatureAngle</i> )                  | The value of this property defines a feature edge. If the surface normal between two adjacent triangles is at least as large as this Feature Angle, a feature edge exists. If Splitting is on, points are duplicated along these feature edges. (See the Splitting property.)  | 30               | The value must be greater than or equal to 0 and less than or equal to 180.   |
| <b>Flip Normals</b><br>( <i>FlipNormals</i> )                    | If the value of this property is 1, this filter will reverse the normal direction (and reorder the points accordingly) for all polygons in the data set; this changes front-facing polygons to back-facing ones, and vice versa. You might want to do this if your viewing position will be inside the data set instead of outside of it.                                  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                 | This property specifies the input to the Normals Generation filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Non-Manifold Traversal</b><br>( <i>NonManifoldTraversal</i> ) | Turn on/off traversal across non-manifold edges. Not traversing non-manifold edges will prevent problems where the consistency of polygonal ordering is corrupted due to topological loops.  | 1                | Only the values 0 and 1 are accepted.   |

|   |  |   |                                       |
|---|--|---|---------------------------------------|
| <b>Piece Invariant</b><br>( <i>PieceInvariant</i> ) | Turn this option to to produce the same results regardless of the number of processors used (i.e., avoid seams along processor boundaries). Turn this off if you do want to process ghost levels and do not mind seams.                                  | 1 | Only the values 0 and 1 are accepted. |
| <b>Splitting</b><br>( <i>Splitting</i> )            | This property controls the splitting of sharp edges. If sharp edges are split (property value = 1), then points are duplicated along these edges, and separate normals are computed for both sets of points to give crisp (rendered) surface definition. | 1 | Only the values 0 and 1 are accepted. |

## Glyph

This filter generates an arrow, cone, cube, cylinder, line, sphere, or 2D glyph at each point of the input data set. The glyphs can be oriented and scaled by point attributes of the input dataset.

The Glyph filter generates a glyph (i.e., an arrow, cone, cube, cylinder, line, sphere, or 2D glyph) at each point in the input dataset. The glyphs can be oriented and scaled by the input point-centered scalars and vectors. The Glyph filter operates on any type of data set. Its output is polygonal. This filter is available on the Toolbar.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Glyph Transform</b><br>( <i>GlyphTransform</i> )                 | The values in this property allow you to specify the transform (translation, rotation, and scaling) to apply to the glyph source.   |                  | The value must be set to one of the following: Transform2.  |
| <b>Input</b><br>( <i>Input</i> )                                    | This property specifies the input to the Glyph filter. This is the dataset to which the glyphs will be applied.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.  |
| <b>Maximum Number of Points</b><br>( <i>MaximumNumberOfPoints</i> ) | The value of this property specifies the maximum number of glyphs that should appear in the output dataset if the value of the UseMaskPoints property is 1. (See the UseMaskPoints property.)       | 5000             | The value must be greater than or equal to 0.   |
| <b>Random Mode</b><br>( <i>RandomMode</i> )                         | If the value of this property is 1, then the points to glyph are chosen randomly. Otherwise the point ids chosen are evenly spaced.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )                     | This property indicates the name of the scalar array on which to operate. The indicated array may be used for scaling the glyphs. (See the SetScaleMode property.)                                  |                  | An array of scalars is required.  |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )                     | This property indicates the name of the vector array on which to operate. The indicated array may be used for scaling and/or orienting the glyphs. (See the SetScaleMode and SetOrient properties.) | 1                | An array of vectors is required.  |
| <b>Orient</b><br>( <i>SetOrient</i> )                               | If this property is set to 1, the glyphs will be oriented based on the selected vector array.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Set Scale Factor</b><br>( <i>SetScaleFactor</i> )                | The value of this property will be used as a multiplier for scaling the glyphs before adding them to the output.  | 1                | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 0.1.<br><br>The value must lie within the range of the selected data array.<br><br>The value must lie within the range of the selected data array. |

|  |  |   |  |
|--|--|---|--|
| <b>Scale Mode</b><br>( <i>SetScaleMode</i> )   | The value of this property specifies how/if the glyphs should be scaled based on the point-centered scalars/vectors in the input dataset.                                | 1 | The value must be one of the following: scalar (0), vector (1), vector_components (2), off (3).  |
| <b>Glyph Type</b><br>( <i>Source</i> )         | This property determines which type of glyph will be placed at the points in the input dataset.  |   | <p>The selected object must be the result of the following: sources (includes readers), glyph_sources.</p> <p>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData.</p> <p>The value must be set to one of the following: ArrowSource, ConeSource, CubeSource, CylinderSource, LineSource, SphereSource, GlyphSource2D.</p> |
| <b>Mask Points</b><br>( <i>UseMaskPoints</i> ) | If the value of this property is set to 1, limit the maximum number of glyphs to the value indicated by MaximumNumberOfPoints. (See the MaximumNumberOfPoints property.) | 1 | Only the values 0 and 1 are accepted.  |

## Glyph With Custom Source

This filter generates a glyph at each point of the input data set. The glyphs can be oriented and scaled by point attributes of the input dataset.

The Glyph filter generates a glyph at each point in the input dataset. The glyphs can be oriented and scaled by the input point-centered scalars and vectors. The Glyph filter operates on any type of data set. Its output is polygonal. This filter is available on the Toolbar.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                                    | This property specifies the input to the Glyph filter. This is the dataset to which the glyphs will be applied.   |                  | <p>The selected object must be the result of the following: sources (includes readers), filters.</p> <p>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.</p> |
| <b>Maximum Number of Points</b><br>( <i>MaximumNumberOfPoints</i> ) | The value of this property specifies the maximum number of glyphs that should appear in the output dataset if the value of the UseMaskPoints property is 1. (See the UseMaskPoints property.)       | 5000             | The value must be greater than or equal to 0.   |
| <b>Random Mode</b><br>( <i>RandomMode</i> )                         | If the value of this property is 1, then the points to glyph are chosen randomly. Otherwise the point ids chosen are evenly spaced.   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )                     | This property indicates the name of the scalar array on which to operate. The indicated array may be used for scaling the glyphs. (See the SetScaleMode property.)                                  |                  | An array of scalars is required.  |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )                     | This property indicates the name of the vector array on which to operate. The indicated array may be used for scaling and/or orienting the glyphs. (See the SetScaleMode and SetOrient properties.) | 1                | An array of vectors is required.  |
| <b>Orient</b><br>( <i>SetOrient</i> )                               | If this property is set to 1, the glyphs will be oriented based on the selected vector array.   | 1                | Only the values 0 and 1 are accepted.   |

|  |  |   |   |
|--|--|---|---|
| <b>Set Scale Factor</b><br>( <i>SetScaleFactor</i> ) | The value of this property will be used as a multiplier for scaling the glyphs before adding them to the output.   | 1 | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 0.1.<br><br>The value must lie within the range of the selected data array.<br><br>The value must lie within the range of the selected data array. |
| <b>Scale Mode</b><br>( <i>SetScaleMode</i> )         | The value of this property specifies how/if the glyphs should be scaled based on the point-centered scalars/vectors in the input dataset.                                | 1 | The value must be one of the following: scalar (0), vector (1), vector_components (2), off (3).   |
| <b>Glyph Type</b><br>( <i>Source</i> )               | This property determines which type of glyph will be placed at the points in the input dataset.  |   | The selected object must be the result of the following: sources (includes readers), glyph_sources.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData.                                     |
| <b>Mask Points</b><br>( <i>UseMaskPoints</i> )       | If the value of this property is set to 1, limit the maximum number of glyphs to the value indicated by MaximumNumberOfPoints. (See the MaximumNumberOfPoints property.) | 1 | Only the values 0 and 1 are accepted.   |

## Gradient

This filter computes gradient vectors for an image/volume.

The Gradient filter computes the gradient vector at each point in an image or volume. This filter uses central differences to compute the gradients. The Gradient filter operates on uniform rectilinear (image) data and produces image data output.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Dimensionality</b><br>( <i>Dimensionality</i> )           | This property indicates whether to compute the gradient in two dimensions or in three. If the gradient is being computed in two dimensions, the X and Y dimensions are used. | 3                | The value must be one of the following: Two (2), Three (3).   |
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Gradient filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |
| <b>Select Input Scalars</b><br>( <i>SelectInputScalars</i> ) | This property lists the name of the array from which to compute the gradient.  |                  | An array of scalars is required.  |

## Gradient Of Unstructured DataSet

Estimate the gradient for each point or cell in any type of dataset.

The Gradient (Unstructured) filter estimates the gradient vector at each point or cell. It operates on any type of `vtkDataSet`, and the output is the same type as the input. If the dataset is a `vtkImageData`, use the Gradient filter instead; it will be more efficient for this type of dataset.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Compute Vorticity</b><br>( <i>ComputeVorticity</i> )       | When this flag is on, the gradient filter will compute the vorticity/curl of a 3 component array.  | 0                | Only the values 0 and 1 are accepted.  |
| <b>Faster Approximation</b><br>( <i>FasterApproximation</i> ) | When this flag is on, the gradient filter will provide a less accurate (but close) algorithm that performs fewer derivative calculations (and is therefore faster). The error contains some smoothing of the output data and some possible errors on the boundary. This parameter has no effect when performing the gradient of cell data. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )                              | This property specifies the input to the Gradient (Unstructured) filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array.<br>The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkPointSet</code> . |
| <b>Result Array Name</b><br>( <i>ResultArrayName</i> )        | This property provides a name for the output array containing the gradient vectors.  | Gradients        |  |
| <b>Scalar Array</b><br>( <i>SelectInputScalars</i> )          | This property lists the name of the scalar array from which to compute the gradient.   |                  | An array of scalars is required.<br>Valid array names will be chosen from point and cell data.   |

## Grid Connectivity

Mass properties of connected fragments for unstructured grids.

This filter works on multiblock unstructured grid inputs and also works in parallel. It ignores any cells with a cell data Status value of 0.

It performs connectivity to distinct fragments separately. It then integrates attributes of the fragments.

| Property                         | Description                                      | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input of the filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkUnstructuredGrid</code> , <code>vtkCompositeDataSet</code> . |

## Group Datasets

Group data sets.

Groups multiple datasets to create a multiblock dataset

| Property                         | Description  | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | This property indicates the inputs to the Group Datasets filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |

## Histogram

Extract a histogram from field data.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Bin Count</b><br>( <i>BinCount</i> )                       | The value of this property specifies the number of bins for the histogram.   | 10               | The value must be greater than or equal to 1 and less than or equal to 256.   |
| <b>Calculate Averages</b><br>( <i>CalculateAverages</i> )     | This option controls whether the algorithm calculates averages of variables other than the primary variable that fall into each bin. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Component</b><br>( <i>Component</i> )                      | The value of this property specifies the array component from which the histogram should be computed.                                | 0                |   |
| <b>Custom Bin Ranges</b><br>( <i>CustomBinRanges</i> )        | Set custom bin ranges to use. These are used only when UseCustomBinRanges is set to true.  | 0 100            | The value must lie within the range of the selected data array.   |
| <b>Input</b><br>( <i>Input</i> )                              | This property specifies the input to the Histogram filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point or cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Select Input Array</b><br>( <i>SelectInputArray</i> )      | This property indicates the name of the array from which to compute the histogram.   |                  | An array of scalars is required.<br><br>Valid array names will be chosen from point and cell data.  |
| <b>Use Custom Bin Ranges</b><br>( <i>UseCustomBinRanges</i> ) | When set to true, CustomBinRanges will be used instead of using the full range for the selected array. By default, set to false.     | 0                | Only the values 0 and 1 are accepted.   |

## Integrate Variables

This filter integrates cell and point attributes.

The Integrate Attributes filter integrates point and cell data over lines and surfaces. It also computes length of lines, area of surface, or volume.

| Property                         | Description   | Default Value(s) | Restrictions   |
|----------------------------------|---|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Integrate Attributes filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Interpolate to Quadrature Points

Create scalar/vector data arrays interpolated to quadrature points.

"Create scalar/vector data arrays interpolated to quadrature points."

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                           | This property specifies the input of the filter.                                  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |
| <b>Select Source Array</b><br>( <i>SelectSourceArray</i> ) | Specifies the offset array from which we interpolate values to quadrature points. |                  | An array of scalars is required.  |

## Intersect Fragments

The Intersect Fragments filter perform geometric intersections on sets of fragments.

The Intersect Fragments filter perform geometric intersections on sets of fragments. The filter takes two inputs, the first containing fragment geometry and the second containing fragment centers. The filter has two outputs. The first is geometry that results from the intersection. The second is a set of points that is an approximation of the center of where each fragment has been intersected.

| Property                                    | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Slice Type</b><br>( <i>CutFunction</i> ) | This property sets the type of intersecting geometry, and associated parameters. |                  | The value must be set to one of the following: Plane, Box, Sphere.   |
| <b>Input</b><br>( <i>Input</i> )            | This input must contain fragment geometry.                                       |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkMultiBlockDataSet. |
| <b>Source</b><br>( <i>Source</i> )          | This input must contain fragment centers.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkMultiBlockDataSet. |

## Iso Volume

This filter extracts cells by clipping cells that have point scalars not in the specified range.

This filter clip away the cells using lower and upper thresholds.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                      | This property specifies the input to the Threshold filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array with 1 components.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Input Scalars</b><br>( <i>SelectInputScalars</i> ) | The value of this property contains the name of the scalar array from which to perform thresholding. |                  | An array of scalars is required.<br>Valid array names will be chosen from point and cell data.  |
| <b>Threshold Range</b><br>( <i>ThresholdBetween</i> ) | The values of this property specify the upper and lower bounds of the thresholding operation.        | 0 0              | The value must lie within the range of the selected data array.   |

## K Means

Compute a statistical model of a dataset and/or assess the dataset with a statistical model.

This filter either computes a statistical model of a dataset or takes such a model as its second input. Then, the model (however it is obtained) may optionally be used to assess the input dataset.

This filter iteratively computes the center of k clusters in a space whose coordinates are specified by the arrays you select. The clusters are chosen as local minima of the sum of square Euclidean distances from each point to its nearest cluster center. The model is then a set of cluster centers. Data is assessed by assigning a cluster center and distance to the cluster to each point in the input data set.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> )    | Specify which type of field data the arrays will be drawn from.   | 0                | Valid array names will be chosen from point and cell data.   |
| <b>Input</b><br>( <i>Input</i> )                     | The input to the filter. Arrays from this dataset will be used for computing statistics and/or assessed by a statistical model. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData, vtkStructuredGrid, vtkPolyData, vtkUnstructuredGrid, vtkTable, vtkGraph. |
| <b>k</b><br>( <i>K</i> )                             | Specify the number of clusters.   | 5                | The value must be greater than or equal to 1.  |
| <b>Max Iterations</b><br>( <i>MaxNumIterations</i> ) | Specify the maximum number of iterations in which cluster centers are moved before the algorithm terminates.                    | 50               | The value must be greater than or equal to 1.  |



|   |   |      |  |
|---|---|------|--|
| <b>Model Input</b><br>( <i>ModelInput</i> )             | A previously-calculated model with which to assess a separate dataset. This input is optional.  |      | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable, vtkMultiBlockDataSet. |
| <b>Variables of Interest</b><br>( <i>SelectArrays</i> ) | Choose arrays whose entries will be used to form observations for statistical analysis.   |      | An array of scalars is required.   |
| <b>Task</b><br>( <i>Task</i> )                          | Specify the task to be performed: modeling and/or assessment.<br><br>1. "Statistics of all the data," creates an output table (or tables) summarizing the <b>entire</b> input dataset;<br><br>2. "Model a subset of the data," creates an output table (or tables) summarizing a <b>randomly-chosen subset</b> of the input dataset;<br><br>3. "Assess the data with a model," adds attributes to the first input dataset using a model provided on the second input port; and<br><br>4. "Model and assess the same data," is really just operations 2 and 3 above applied to the same input dataset. The model is first trained using a fraction of the input data and then the entire dataset is assessed using that model.<br><br>When the task includes creating a model (i.e., tasks 2, and 4), you may adjust the fraction of the input dataset used for training. You should avoid using a large fraction of the input data for training as you will then not be able to detect overfitting. The <i>Training fraction</i> setting will be ignored for tasks 1 and 3. | 3    | The value must be one of the following:<br>Statistics of all the data (0), Model a subset of the data (1), Assess the data with a model (2), Model and assess the same data (3).   |
| <b>Tolerance</b><br>( <i>Tolerance</i> )                | Specify the relative tolerance that will cause early termination.   | 0.01 | The value must be greater than or equal to 0 and less than or equal to 1.  |
| <b>Training Fraction</b><br>( <i>TrainingFraction</i> ) | Specify the fraction of values from the input dataset to be used for model fitting. The exact set of values is chosen at random from the dataset.   | 0.1  | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Level Scalars

The Level Scalars filter uses colors to show levels of a hierarchical dataset.

The Level Scalars filter uses colors to show levels of a hierarchical dataset.

| Property                         | Description  | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Level Scalars filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkHierarchicalBoxDataSet. |

## Linear Extrusion

This filter creates a swept surface defined by translating the input along a vector.

The Linear Extrusion filter creates a swept surface by translating the input dataset along a specified vector. This filter is intended to operate on 2D polygonal data. This filter operates on polygonal data and produces polygonal data output.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Capping</b><br>( <i>Capping</i> )                | The value of this property indicates whether to cap the ends of the swept surface. Capping works by placing a copy of the input dataset on either end of the swept surface, so it behaves properly if the input is a 2D surface composed of filled polygons. If the input dataset is a closed solid (e.g., a sphere), then if capping is on (i.e., this property is set to 1), two copies of the data set will be displayed on output (the second translated from the first one along the specified vector). If instead capping is off (i.e., this property is set to 0), then an input closed solid will produce no output. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                    | This property specifies the input to the Linear Extrusion filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Piece Invariant</b><br>( <i>PieceInvariant</i> ) | The value of this property determines whether the output will be the same regardless of the number of processors used to compute the result. The difference is whether there are internal polygonal faces on the processor boundaries. A value of 1 will keep the results the same; a value of 0 will allow internal faces on processor boundaries.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Scale Factor</b><br>( <i>ScaleFactor</i> )       | The value of this property determines the distance along the vector the dataset will be translated. (A scale factor of 0.5 will move the dataset half the length of the vector, and a scale factor of 2 will move it twice the vector's length.)   | 1                |   |
| <b>Vector</b><br>( <i>Vector</i> )                  | The value of this property indicates the X, Y, and Z components of the vector along which to sweep the input dataset.  | 0 0 1            |   |

## Loop Subdivision

This filter iteratively divides each triangle into four triangles. New points are placed so the output surface is smooth.

The Loop Subdivision filter increases the granularity of a polygonal mesh. It works by dividing each triangle in the input into four new triangles. It is named for Charles Loop, the person who devised this subdivision scheme. This filter only operates on triangles, so a data set that contains other types of polygons should be passed through the Triangulate filter before applying this filter to it. This filter only operates on polygonal data (specifically triangle meshes), and it produces polygonal output.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                                 | This property specifies the input to the Loop Subdivision filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Number of Subdivisions</b><br>( <i>NumberOfSubdivisions</i> ) | Set the number of subdivision iterations to perform. Each subdivision divides single triangles into four new triangles. | 1                | The value must be greater than or equal to 1 and less than or equal to 4.   |

## Mask Points

Reduce the number of points. This filter is often used before glyphing. Generating vertices is an option.

The Mask Points filter reduces the number of points in the dataset. It operates on any type of dataset, but produces only points / vertices as output. This filter is often used before the Glyph filter, but the basic point-masking functionality is also available on the Properties page for the Glyph filter.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Generate Vertices</b><br>( <i>GenerateVertices</i> )             | This property specifies whether to generate vertex cells as the topography of the output. If set to 1, the geometry (vertices) will be displayed in the rendering window; otherwise no geometry will be displayed.  | 0                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )                                    | This property specifies the input to the Mask Points filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Maximum Number of Points</b><br>( <i>MaximumNumberOfPoints</i> ) | The value of this property indicates the maximum number of points in the output dataset.  | 5000             | The value must be greater than or equal to 0.  |
| <b>Offset</b><br>( <i>Offset</i> )                                  | The value of this property indicates the point in the input dataset from which to start masking.  | 0                | The value must be greater than or equal to 0.  |
| <b>On Ratio</b><br>( <i>OnRatio</i> )                               | The value of this property specifies the ratio of points to retain in the output. (For example, if the on ratio is 3, then the output will contain 1/3 as many points -- up to the value of the MaximumNumberOfPoints property -- as the input.)                            | 2                | The value must be greater than or equal to 1.  |
| <b>Random</b><br>( <i>RandomMode</i> )                              | If the value of this property is set to 0, then the points in the output will be randomly selected from the input; otherwise this filter will subsample regularly. Selecting points at random is helpful to avoid striping when masking the points of a structured dataset. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Single Vertex Per Cell</b><br>( <i>SingleVertexPerCell</i> )     | Tell filter to only generate one vertex per cell instead of multiple vertices in one cell.  | 0                | Only the values 0 and 1 are accepted.  |

## Material Interface Filter

The Material Interface filter finds volumes in the input data containing material above a certain material fraction.

The Material Interface filter finds voxels inside of which a material fraction (or normalized amount of material) is higher than a given threshold. As these voxels are identified surfaces enclosing adjacent voxels above the threshold are generated. The resulting volume and its surface are what we call a fragment. The filter has the ability to compute various volumetric attributes such as fragment volume, mass, center of mass as well as volume and mass weighted averages for any of the fields present. Any field selected for such computation will be also be copied into the fragment surface's point data for visualization. The filter also has the ability to generate Oriented Bounding Boxes (OBB) for each fragment.

The data generated by the filter is organized in three outputs. The "geometry" output, containing the fragment surfaces. The "statistics" output, containing a point set of the centers of mass. The "obb representation" output, containing OBB representations (poly data). All computed attributes are copied into the statistics and geometry output. The obb representation output is used for validation and debugging purposes and is turned off by default.

To measure the size of craters, the filter can invert a volume fraction and clip the volume fraction with a sphere and/or a plane.

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Clip Type</b><br>( <i>ClipFunction</i> )                                | This property sets the type of clip geometry, and associated parameters.   |                  | The value must be set to one of the following: None, Plane, Sphere.  |
| <b>Compute OBB</b><br>( <i>ComputeOBB</i> )                                | Compute Object Oriented Bounding boxes (OBB). When active the result of this computation is copied into the statistics output. In the case that the filter is built in its validation mode, the OBB's are rendered.                            | 0                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )   | Input to the filter can be a hierarchical box data set containing image data or a multi-block of rectilinear grids.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkHierarchicalBoxDataSet. |
| <b>Invert Volume Fraction</b><br>( <i>InvertVolumeFraction</i> )           | Inverting the volume fraction generates the negative of the material. It is useful for analyzing craters.  | 0                | Only the values 0 and 1 are accepted.  |
| <b>Material Fraction Threshold</b><br>( <i>MaterialFractionThreshold</i> ) | Material fraction is defined as normalized amount of material per voxel. Any voxel in the input data set with a material fraction greater than this value is included in the output data set.  | 0.5              | The value must be greater than or equal to 0.08 and less than or equal to 1.   |
| <b>Output Base Name</b><br>( <i>OutputBaseName</i> )                       | This property specifies the base including path of where to write the statistics and geometry output text files. It follows the pattern "/path/to/folder/and/file" here file has no extension, as the filter will generate a unique extension. |                  |  |

|  |   |   |                                       |
|--|---|---|---------------------------------------|
| <b>Select Mass Arrays</b><br>( <i>SelectMassArray</i> )                            | Mass arrays are paired with material fraction arrays. This means that the first selected material fraction array is paired with the first selected mass array, and so on sequentially. As the filter identifies voxels meeting the minimum material fraction threshold, these voxel's mass will be used in fragment center of mass and mass calculation.<br><br>A warning is generated if no mass array is selected for an individual material fraction array. However, in that case the filter will run without issue because the statistics output can be generated using fragments' centers computed from axis aligned bounding boxes. |   | An array of scalars is required.      |
| <b>Compute mass weighted average over:</b><br>( <i>SelectMassWtdAvgArray</i> )     | For arrays selected a mass weighted average is computed. These arrays are also copied into fragment geometry cell data as the fragment surfaces are generated.  |   | An array of scalars is required.      |
| <b>Select Material Fraction Arrays</b><br>( <i>SelectMaterialArray</i> )           | Material fraction is defined as normalized amount of material per voxel. It is expected that arrays containing material fraction data has been down converted to a unsigned char.   |   | An array of scalars is required.      |
| <b>Compute volume weighted average over:</b><br>( <i>SelectVolumeWtdAvgArray</i> ) | For arrays selected a volume weighted average is computed. The values of these arrays are also copied into fragment geometry cell data as the fragment surfaces are generated.  |   | An array of scalars is required.      |
| <b>Write Geometry Output</b><br>( <i>WriteGeometryOutput</i> )                     | If this property is set, then the geometry output is written to a text file. The file name will be coconstructed using the path in the "Output Base Name" widget.   | 0 | Only the values 0 and 1 are accepted. |
| <b>Write Statistics Output</b><br>( <i>WriteStatisticsOutput</i> )                 | If this property is set, then the statistics output is written to a text file. The file name will be coconstructed using the path in the "Output Base Name" widget.   | 0 | Only the values 0 and 1 are accepted. |

## Median

Compute the median scalar values in a specified neighborhood for image/volume datasets.

The Median filter operates on uniform rectilinear (image or volume) data and produces uniform rectilinear output. It replaces the scalar value at each pixel / voxel with the median scalar value in the specified surrounding neighborhood. Since the median operation removes outliers, this filter is useful for removing high-intensity, low-probability noise (shot noise).

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Median filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |
| <b>Kernel Size</b><br>( <i>KernelSize</i> )                  | The value of this property specifies the number of pixels/voxels in each dimension to use in computing the median to assign to each pixel/voxel. If the kernel size in a particular dimension is 1, then the median will not be computed in that direction. | 1 1 1            |   |
| <b>Select Input Scalars</b><br>( <i>SelectInputScalars</i> ) | The value of this property lists the name of the scalar array to use in computing the median.   |                  | An array of scalars is required.  |

## Merge Blocks

`vtkCompositeDataToUnstructuredGridFilter` appends all `vtkDataSet` leaves of the input composite dataset to a single unstructure grid. The subtree to be combined can be chosen using the `SubTreeCompositeIndex`. If the `SubTreeCompositeIndex` is a leaf node, then no appending is required.

| Property                         | Description                      | Default Value(s) | Restrictions   |
|----------------------------------|----------------------------------|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | Set the input composite dataset. |                  | The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkCompositeDataSet</code> . |

## Mesh Quality

This filter creates a new cell array containing a geometric measure of each cell's fitness. Different quality measures can be chosen for different cell shapes.

This filter creates a new cell array containing a geometric measure of each cell's fitness. Different quality measures can be chosen for different cell shapes. Supported shapes include triangles, quadrilaterals, tetrahedra, and hexahedra. For other shapes, a value of 0 is assigned.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Hex Quality Measure</b><br>( <i>HexQualityMeasure</i> )   | This property indicates which quality measure will be used to evaluate hexahedral quality.    | 5                | The value must be one of the following: Diagonal (21), Dimension (22), Distortion (15), Edge Ratio (0), Jacobian (25), Maximum Edge Ratio (16), Maximum Aspect Frobenius (5), Mean Aspect Frobenius (4), Oddy (23), Relative Size Squared (12), Scaled Jacobian (10), Shape (13), Shape and Size (14), Shear (11), Shear and Size (24), Skew (17), Stretch (20), Taper (18), Volume (19).  |
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Mesh Quality filter.                                 |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkDataSet</code> .   |
| <b>Quad Quality Measure</b><br>( <i>QuadQualityMeasure</i> ) | This property indicates which quality measure will be used to evaluate quadrilateral quality. | 0                | The value must be one of the following: Area (28), Aspect Ratio (1), Condition (9), Distortion (15), Edge Ratio (0), Jacobian (25), Maximum Aspect Frobenius (5), Maximum Aspect Frobenius (5), Maximum Edge Ratio (16), Mean Aspect Frobenius (4), Minimum Angle (6), Oddy (23), Radius Ratio (2), Relative Size Squared (12), Scaled Jacobian (10), Shape (13), Shape and Size (14), Shear (11), Shear and Size (24), Skew (17), Stretch (20), Taper (18), Warpage (26). |

|  |  |   |   |
|--|--|---|---|
| <b>Tet Quality Measure</b><br>( <i>TetQualityMeasure</i> )           | This property indicates which quality measure will be used to evaluate tetrahedral quality. The radius ratio is the size of a sphere circumscribed by a tetrahedron's 4 vertices divided by the size of a circle tangent to a tetrahedron's 4 faces. The edge ratio is the ratio of the longest edge length to the shortest edge length. The collapse ratio is the minimum ratio of height of a vertex above the triangle opposite it divided by the longest edge of the opposing triangle across all vertex/triangle pairs. | 2 | The value must be one of the following: Edge Ratio (0), Aspect Beta (29), Aspect Gamma (27), Aspect Frobenius (3), Aspect Ratio (1), Collapse Ratio (7), Condition (9), Distortion (15), Jacobian (25), Minimum Dihedral Angle (6), Radius Ratio (2), Relative Size Squared (12), Scaled Jacobian (10), Shape (13), Shape and Size (14), Volume (19). |
| <b>Triangle Quality Measure</b><br>( <i>TriangleQualityMeasure</i> ) | This property indicates which quality measure will be used to evaluate triangle quality. The radius ratio is the size of a circle circumscribed by a triangle's 3 vertices divided by the size of a circle tangent to a triangle's 3 edges. The edge ratio is the ratio of the longest edge length to the shortest edge length.  | 2 | The value must be one of the following: Area (28), Aspect Ratio (1), Aspect Frobenius (3), Condition (9), Distortion (15), Edge Ratio (0), Maximum Angle (8), Minimum Angle (6), Scaled Jacobian (10), Radius Ratio (2), Relative Size Squared (12), Shape (13), Shape and Size (14).   |

## Multicorrelative Statistics

Compute a statistical model of a dataset and/or assess the dataset with a statistical model.

This filter either computes a statistical model of a dataset or takes such a model as its second input. Then, the model (however it is obtained) may optionally be used to assess the input dataset.

This filter computes the covariance matrix for all the arrays you select plus the mean of each array. The model is thus a multivariate Gaussian distribution with the mean vector and variances provided. Data is assessed using this model by computing the Mahalanobis distance for each input point. This distance will always be positive.

The learned model output format is rather dense and can be confusing, so it is discussed here. The first filter output is a multiblock dataset consisting of 2 tables:

1. Raw covariance data.
2. Covariance matrix and its Cholesky decomposition.

The raw covariance table has 3 meaningful columns: 2 titled "Column1" and "Column2" whose entries generally refer to the N arrays you selected when preparing the filter and 1 column titled "Entries" that contains numeric values. The first row will always contain the number of observations in the statistical analysis. The next N rows contain the mean for each of the N arrays you selected. The remaining rows contain covariances of pairs of arrays.

The second table (covariance matrix and Cholesky decomposition) contains information derived from the raw covariance data of the first table. The first N rows of the first column contain the name of one array you selected for analysis. These rows are followed by a single entry labeled "Cholesky" for a total of N+1 rows. The second column, Mean contains the mean of each variable in the first N entries and the number of observations processed in the final (N+1) row.

The remaining columns (there are N, one for each array) contain 2 matrices in triangular format. The upper right triangle contains the covariance matrix (which is symmetric, so its lower triangle may be inferred). The lower left triangle contains the Cholesky decomposition of the covariance matrix (which is triangular, so its upper triangle is zero). Because the diagonal must be stored for both matrices, an additional row is required  $\sqrt{N+1}$  hence the N+1 rows and the final entry of the column named "Column".

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> )       | Specify which type of field data the arrays will be drawn from.   | 0                | Valid array names will be chosen from point and cell data.   |
| <b>Input</b><br>( <i>Input</i> )                        | The input to the filter. Arrays from this dataset will be used for computing statistics and/or assessed by a statistical model.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The dataset must contain a point or cell array.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData, vtkStructuredGrid, vtkPolyData, vtkUnstructuredGrid, vtkTable, vtkGraph. |
| <b>Model Input</b><br>( <i>ModelInput</i> )             | A previously-calculated model with which to assess a separate dataset. This input is optional.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable, vtkMultiBlockDataSet.   |
| <b>Variables of Interest</b><br>( <i>SelectArrays</i> ) | Choose arrays whose entries will be used to form observations for statistical analysis.   |                  | An array of scalars is required.   |
| <b>Task</b><br>( <i>Task</i> )                          | Specify the task to be performed: modeling and/or assessment.<br><ol style="list-style-type: none"> <li>1. "Statistics of all the data," creates an output table (or tables) summarizing the <b>entire</b> input dataset;</li> <li>2. "Model a subset of the data," creates an output table (or tables) summarizing a <b>randomly-chosen subset</b> of the input dataset;</li> <li>3. "Assess the data with a model," adds attributes to the first input dataset using a model provided on the second input port; and</li> <li>4. "Model and assess the same data," is really just operations 2 and 3 above applied to the same input dataset. The model is first trained using a fraction of the input data and then the entire dataset is assessed using that model.</li> </ol> <p>When the task includes creating a model (i.e., tasks 2, and 4), you may adjust the fraction of the input dataset used for training. You should avoid using a large fraction of the input data for training as you will then not be able to detect overfitting. The <i>Training fraction</i> setting will be ignored for tasks 1 and 3.</p> | 3                | The value must be one of the following:<br>Statistics of all the data (0), Model a subset of the data (1), Assess the data with a model (2), Model and assess the same data (3).   |
| <b>Training Fraction</b><br>( <i>TrainingFraction</i> ) | Specify the fraction of values from the input dataset to be used for model fitting. The exact set of values is chosen at random from the dataset.   | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Normal Glyphs

Filter computing surface normals.

Filter computing surface normals.



| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Consistency</b><br>( <i>Consistency</i> )                    | The value of this property controls whether consistent polygon ordering is enforced. Generally the normals for a data set should either all point inward or all point outward. If the value of this property is 1, then this filter will reorder the points of cells that whose normal vectors are oriented the opposite direction from the rest of those in the data set. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Maximum Number of Points</b><br>( <i>Glyph Max. Points</i> ) | The value of this property specifies the maximum number of glyphs that should appear in the output dataset if the value of the UseMaskPoints property is 1. (See the UseMaskPoints property.)  | 5000             | The value must be greater than or equal to 0.   |
| <b>Random Mode</b><br>( <i>Glyph Random Mode</i> )              | If the value of this property is 1, then the points to glyph are chosen randomly. Otherwise the point ids chosen are evenly spaced.  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Set Scale Factor</b><br>( <i>Glyph Scale Factor</i> )        | The value of this property will be used as a multiplier for scaling the glyphs before adding them to the output.   | 1                | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 0.1.<br><br>The value must lie within the range of the selected data array.<br><br>The value must lie within the range of the selected data array. |
| <b>Input</b><br>( <i>Input</i> )                                | This property specifies the input to the Extract Surface filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.  |
| <b>Invert</b><br>( <i>InvertArrow</i> )                         | Inverts the arrow direction.   | 0                | Only the values 0 and 1 are accepted.   |

## Octree Depth Limit

This filter takes in a octree and produces a new octree which is no deeper than the maximum specified depth level.

The Octree Depth Limit filter takes in an octree and produces a new octree that is nowhere deeper than the maximum specified depth level. The attribute data of pruned leaf cells are integrated in to their ancestors at the cut level.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Octree Depth Limit filter.          |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkHyperOctree. |
| <b>Maximum Level</b><br>( <i>MaximumLevel</i> ) | The value of this property specifies the maximum depth of the output octree. | 4                | The value must be greater than or equal to 3 and less than or equal to 255.  |

## Octree Depth Scalars

This filter adds a scalar to each leaf of the octree that represents the leaf's depth within the tree.

The vtkHyperOctreeDepth filter adds a scalar to each leaf of the octree that represents the leaf's depth within the tree.

| Property                         | Description   | Default Value(s) | Restrictions   |
|----------------------------------|---|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Octree Depth Scalars filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkHyperOctree. |

## Outline

This filter generates a bounding box representation of the input.

The Outline filter generates an axis-aligned bounding box for the input dataset. This filter operates on any type of dataset and produces polygonal output.

| Property                         | Description  | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Outline filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Outline Corners

This filter generates a bounding box representation of the input. It only displays the corners of the bounding box.

The Outline Corners filter generates the corners of a bounding box for the input dataset. This filter operates on any type of dataset and produces polygonal output.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Corner Factor</b><br>( <i>CornerFactor</i> ) | The value of this property sets the size of the corners as a percentage of the length of the corresponding bounding box edge. | 0.2              | The value must be greater than or equal to 0.001 and less than or equal to 0.5.  |
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Outline Corners filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Outline Curvilinear DataSet

This filter generates an outline representation of the input.

The Outline filter generates an outline of the outside edges of the input dataset, rather than the dataset's bounding box. This filter operates on structured grid datasets and produces polygonal output.

| Property                         | Description  | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the outline (curvilinear) filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkStructuredGrid. |

## Particle Pathlines

Creates polylines representing pathlines of animating particles

Particle Pathlines takes any dataset as input, it extracts the point locations of all cells over time to build up a polyline trail. The point number (index) is used as the 'key' if the points are randomly changing their respective order in the points list, then you should specify a scalar that represents the unique ID. This is intended to handle the output of a filter such as the TemporalStreamTracer.

| Property   | Description  | Default Value(s)    | Restrictions  |
|--|--|---------------------|---|
| <b>Id Channel Array</b><br>( <i>IdChannelArray</i> ) | Specify the name of a scalar array which will be used to fetch the index of each point. This is necessary only if the particles change position (Id order) on each time step. The Id can be used to identify particles at each step and hence track them properly. If this array is set to "Global or Local IDs", the global point ids are used if they exist or the point index is otherwise. | Global or Local IDs | An array of scalars is required.  |
| <b>Input</b><br>( <i>Input</i> )                     | The input cells to create pathlines for.   |                     | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |

|   |  |       |  |
|---|--|-------|--|
| <b>Mask Points</b><br>(MaskPoints)            | Set the number of particles to track as a ratio of the input. Example: setting MaskPoints to 10 will track every 10th point.   | 100   |  |
| <b>Max Step Distance</b><br>(MaxStepDistance) | If a particle disappears from one end of a simulation and reappears on the other side, the track left will be unrepresentative. Set a MaxStepDistance{x,y,z} which acts as a threshold above which if a step occurs larger than the value (for the dimension), the track will be dropped and restarted after the step. (ie the part before the wrap around will be dropped and the newer part kept). | 1 1 1 |  |
| <b>Max Track Length</b><br>(MaxTrackLength)   | If the Particles being traced animate for a long time, the trails or traces will become long and stringy. Setting the MaxTraceTimeLength will limit how much of the trace is displayed. Tracks longer then the Max will disappear and the trace will appear like a snake of fixed length which progresses as the particle moves. This length is given with respect to timesteps.                     | 25    |  |
| <b>Selection</b><br>(Selection)               | Set a second input, which is a selection. Particles with the same Id in the selection as the primary input will be chosen for pathlines Note that you must have the same IdChannelArray in the selection as the input  |       | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## ParticleTracer

Trace Particles through time in a vector field.

The Particle Trace filter generates pathlines in a vector field from a collection of seed points. The vector field used is selected from the Vectors menu, so the input data set is required to have point-centered vectors. The Seed portion of the interface allows you to select whether the seed points for this integration lie in a point cloud or along a line. Depending on which is selected, the appropriate 3D widget (point or line widget) is displayed along with traditional user interface controls for positioning the point cloud or line within the data set. Instructions for using the 3D widgets and the corresponding manual controls can be found in section 7.4.

This filter operates on any type of data set, provided it has point-centered vectors. The output is polygonal data containing polylines. This filter is available on the Toolbar.

| Property   | Description   | Default Value(s) | Restrictions                          |
|--|---|------------------|---------------------------------------|
| <b>Compute Vorticity</b><br>(ComputeVorticity)                         | Compute vorticity and angular rotation of particles as they progress  | 1                | Only the values 0 and 1 are accepted. |
| <b>Enable Particle Writing</b><br>(EnableParticleWriting)              | Turn On/Off particle writing  | 0                | Only the values 0 and 1 are accepted. |
| <b>Force ReInjection Every NSteps</b><br>(ForceReInjectionEveryNSteps) | When animating particles, it is nice to inject new ones every Nth step to produce a continuous flow. Setting ForceReInjectionEveryNSteps to a non zero value will cause the particle source to reinject particles every Nth step even if it is otherwise unchanged. Note that if the particle source is also animated, this flag will be redundant as the particles will be reinjected whenever the source changes anyway | 1                |                                       |
| <b>Ignore Pipeline Time</b><br>(IgnorePipelineTime)                    | Ignore the TIME_ requests made by the pipeline and only use the TimeStep set manually   | 0                | Only the values 0 and 1 are accepted. |

|  |   |                                      |  |
|--|---|--------------------------------------|--|
| <b>Initial Integration Step</b><br>( <i>InitialIntegrationStep</i> ) | Specify the Initial step size used for line integration, expressed in: LENGTH_UNIT = 1<br>CELL_LENGTH_UNIT = 2 (either the starting size for an adaptive integrator, e.g., RK45, or the constant / fixed size for non-adaptive ones, i.e., RK2 and RK4) | 0.25                                 |  |
| <b>Input</b><br>( <i>Input</i> )                                     | Specify which is the Input of the StreamTracer filter.  |                                      | The selected object must be the result of the following:<br>sources (includes readers), filters.<br>The dataset must contain a point array with 3 components.<br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkDataObject. |
| <b>Particle File Name</b><br>( <i>ParticleFileName</i> )             | Provide a name for the particle file generated if writing is enabled  | /project/cvis/biddisco/ptracer/run-1 |  |
| <b>Select Input Vectors</b><br>( <i>SelectInputVectors</i> )         | Specify which vector array should be used for the integration through that filter.  |                                      | An array of vectors is required.   |
| <b>Source</b><br>( <i>Source</i> )                                   | Specify the seed dataset. Typically from where the vector field integration should begin. Usually a point/radius or a line with a given resolution.   |                                      | The selected object must be the result of the following:<br>sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.  |
| <b>Static Mesh</b><br>( <i>StaticMesh</i> )                          | Force the use of static mesh optimizations  | 0                                    | Only the values 0 and 1 are accepted.  |
| <b>Static Seeds</b><br>( <i>StaticSeeds</i> )                        | Force the use of static seed optimizations  | 1                                    | Only the values 0 and 1 are accepted.  |
| <b>Term. Speed</b><br>( <i>TerminalSpeed</i> )                       | If at any point the speed is below the value of this property, the integration is terminated.   | 1e-12                                |  |
| <b>Termination Time</b><br>( <i>TerminationTime</i> )                | Setting TerminationTime to a positive value will cause particles to terminate when the time is reached. Use a value of zero to disable termination. The units of time should be consistent with the primary time variable.                              | 0                                    |  |
| <b>Termination Time Unit</b><br>( <i>TerminationTimeUnit</i> )       | The termination time may be specified as TimeSteps or Simulation time   | 1                                    | The value must be one of the following:<br>Simulation Time (0), TimeSteps (1).   |
| <b>Time Step</b><br>( <i>TimeStep</i> )                              | Set/Get the TimeStep. This is the primary means of advancing the particles. The TimeStep should be animated and this will drive the pipeline forcing timesteps to be fetched from upstream.   | 0                                    |  |

## Plot Data

This filter prepare arbitrary data to be plotted in any of the plots.

By default the data is shown in a XY line plot.

| Property                         | Description | Default Value(s) | Restrictions  |
|----------------------------------|-------------|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | The input.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |

## Plot Global Variables Over Time

Extracts and plots data in field data over time.

This filter extracts the variables that reside in a dataset's field data and are defined over time. The output is a 1D rectilinear grid where the x coordinates correspond to time (the same array is also copied to a point array named Time or TimeData (if Time exists in the input)).

| Property                         | Description                                      | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | The input from which the selection is extracted. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Plot On Intersection Curves

Extracts the edges in a 2D plane and plots them

Extracts the surface, intersect it with a 2D plane.

Plot the resulting polylines.

| Property                                   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )           | This property specifies the input to the Extract Surface filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Slice Type</b><br>( <i>Slice Type</i> ) | This property sets the parameters of the slice function.         |                  | The value must be set to one of the following: Plane, Box, Sphere.   |

## Plot On Sorted Lines

| Property | Description | Default Value(s) | Restrictions |
|----------|-------------|------------------|--------------|
|----------|-------------|------------------|--------------|

|                                  |   |  |   |
|----------------------------------|---|--|---|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Plot Edges filter. |  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
|----------------------------------|---|--|---|

## Plot Over Line

Sample data attributes at the points along a line. Probed lines will be displayed in a graph of the attributes.

The Plot Over Line filter samples the data set attributes of the current data set at the points along a line. The values of the point-centered variables along that line will be displayed in an XY Plot. This filter uses interpolation to determine the values at the selected point, whether or not it lies at an input point. The Probe filter operates on any type of data and produces polygonal output (a line).

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                           | This property specifies the dataset from which to obtain probe values.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point or cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkDataSet,<br>vtkCompositeDataSet. |
| <b>Pass Partial Arrays</b><br>( <i>PassPartialArrays</i> ) | When dealing with composite datasets, partial arrays are common i.e. data-arrays that are not available in all of the blocks. By default, this filter only passes those point and cell data-arrays that are available in all the blocks i.e. partial array are removed. When PassPartialArrays is turned on, this behavior is changed to take a union of all arrays present thus partial arrays are passed as well. However, for composite dataset input, this filter still produces a non-composite output. For all those locations in a block of where a particular data array is missing, this filter uses vtkMath::Nan() for double and float arrays, while 0 for all other types of arrays i.e int, char etc. | 1                | Only the values 0 and 1 are accepted.  |
| <b>Probe Type</b><br>( <i>Source</i> )                     | This property specifies the dataset whose geometry will be used in determining positions to probe.   |                  | The selected object must be the result of the following: sources (includes readers).<br><br>The value must be set to one of the following:<br>HighResLineSource.   |

## Plot Selection Over Time

Extracts selection over time and then plots it.

This filter extracts the selection over time, i.e. cell and/or point variables at a cells/point selected are extracted over time

The output multi-block consists of 1D rectilinear grids where the x coordinate corresponds to time (the same array is also copied to a point array named Time or TimeData (if Time exists in the input)).

If selection input is a Location based selection then the point values are interpolated from the nearby cells, ie those of the cell the location lies in.

| Property                                 | Description                                      | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )         | The input from which the selection is extracted. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet, vtkTable, vtkCompositeDataSet. |
| <b>Selection</b><br>( <i>Selection</i> ) | The input that provides the selection object.    |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkSelection.                              |

## Point Data to Cell Data

Create cell attributes by averaging point attributes.

The Point Data to Cell Data filter averages the values of the point attributes of the points of a cell to compute cell attributes. This filter operates on any type of dataset, and the output dataset is the same type as the input.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                   | This property specifies the input to the Point Data to Cell Data filter.  |                  | Once set, the input dataset type cannot be changed.<br><br>The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Pass Point Data</b><br>( <i>PassPointData</i> ) | The value of this property controls whether the input point data will be passed to the output. If set to 1, then the input point data is passed through to the output; otherwise, only generated cell data is placed into the output. | 0                | Only the values 0 and 1 are accepted.  |

## Principal Component Analysis

Compute a statistical model of a dataset and/or assess the dataset with a statistical model.

This filter either computes a statistical model of a dataset or takes such a model as its second input. Then, the model (however it is obtained) may optionally be used to assess the input dataset.

This filter performs additional analysis above and beyond the multicorrelative filter. It computes the eigenvalues and eigenvectors of the covariance matrix from the multicorrelative filter. Data is then assessed by projecting the original tuples into a possibly lower-dimensional space.

Since the PCA filter uses the multicorrelative filter's analysis, it shares the same raw covariance table specified in the multicorrelative documentation. The second table in the multiblock dataset comprising the model output is an expanded version of the multicorrelative version.

As with the multicorrelative filter, the second model table contains the mean values, the upper-triangular portion of the symmetric covariance matrix, and the non-zero lower-triangular portion of the Cholesky decomposition of the



covariance matrix. Below these entries are the eigenvalues of the covariance matrix (in the column labeled "Mean") and the eigenvectors (as row vectors) in an additional NxN matrix.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Attribute Mode</b><br>( <i>AttributeMode</i> ) | Specify which type of field data the arrays will be drawn from.  | 0                | Valid array names will be chosen from point and cell data.  |
| <b>Basis Energy</b><br>( <i>BasisEnergy</i> )     | The minimum energy to use when determining the dimensionality of the new space into which the assessment will project tuples.  | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.   |
| <b>Basis Scheme</b><br>( <i>BasisScheme</i> )     | <p>When reporting assessments, should the full eigenvector decomposition be used to project the original vector into the new space (Full basis), or should a fixed subset of the decomposition be used (Fixed-size basis), or should the projection be clipped to preserve at least some fixed "energy" (Fixed-energy basis)?</p> <p>As an example, suppose the variables of interest were {A,B,C,D,E} and that the eigenvalues of the covariance matrix for these were {5,2,1.5,1,.5}. If the "Full basis" scheme is used, then all 5 components of the eigenvectors will be used to project each {A,B,C,D,E}-tuple in the original data into a new 5-components space.</p> <p>If the "Fixed-size" scheme is used and the "Basis Size" property is set to 4, then only the first 4 eigenvector components will be used to project each {A,B,C,D,E}-tuple into the new space and that space will be of dimension 4, not 5.</p> <p>If the "Fixed-energy basis" scheme is used and the "Basis Energy" property is set to 0.8, then only the first 3 eigenvector components will be used to project each {A,B,C,D,E}-tuple into the new space, which will be of dimension 3. The number 3 is chosen because 3 is the lowest N for which the sum of the first N eigenvalues divided by the sum of all eigenvalues is larger than the specified "Basis Energy" (i.e., <math>(5+2+1.5)/10 = 0.85 &gt; 0.8</math>).</p> | 0                | The value must be one of the following: Full basis (0), Fixed-size basis (1), Fixed-energy basis (2).   |
| <b>Basis Size</b><br>( <i>BasisSize</i> )         | The maximum number of eigenvector components to use when projecting into the new space.  | 2                | The value must be greater than or equal to 1.   |
| <b>Input</b><br>( <i>Input</i> )                  | The input to the filter. Arrays from this dataset will be used for computing statistics and/or assessed by a statistical model.  |                  | <p>The selected object must be the result of the following: sources (includes readers), filters.</p> <p>The dataset must contain a point or cell array.</p> <p>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData, vtkStructuredGrid, vtkPolyData, vtkUnstructuredGrid, vtkTable, vtkGraph.</p> |
| <b>Model Input</b><br>( <i>ModelInput</i> )       | A previously-calculated model with which to assess a separate dataset. This input is optional.   |                  | <p>The selected object must be the result of the following: sources (includes readers), filters.</p> <p>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable, vtkMultiBlockDataSet.</p>   |

|   |   |     |  |
|---|---|-----|--|
| <b>Normalization Scheme</b><br>( <i>NormalizationScheme</i> ) | Before the eigenvector decomposition of the covariance matrix takes place, you may normalize each (i,j) entry by $\sqrt{\text{cov}(i,i) * \text{cov}(j,j)}$ . This implies that the variance of each variable of interest should be of equal importance.  | 2   | The value must be one of the following:<br>No normalization (0), Normalize using covariances (3).  |
| <b>Variables of Interest</b><br>( <i>SelectArrays</i> )       | Choose arrays whose entries will be used to form observations for statistical analysis.   |     | An array of scalars is required.   |
| <b>Task</b><br>( <i>Task</i> )                                | Specify the task to be performed: modeling and/or assessment.<br><br>1. "Statistics of all the data," creates an output table (or tables) summarizing the <b>entire</b> input dataset;<br>2. "Model a subset of the data," creates an output table (or tables) summarizing a <b>randomly-chosen subset</b> of the input dataset;<br>3. "Assess the data with a model," adds attributes to the first input dataset using a model provided on the second input port; and<br>4. "Model and assess the same data," is really just operations 2 and 3 above applied to the same input dataset. The model is first trained using a fraction of the input data and then the entire dataset is assessed using that model.<br><br>When the task includes creating a model (i.e., tasks 2, and 4), you may adjust the fraction of the input dataset used for training. You should avoid using a large fraction of the input data for training as you will then not be able to detect overfitting. The <i>Training fraction</i> setting will be ignored for tasks 1 and 3. | 3   | The value must be one of the following:<br>Statistics of all the data (0), Model a subset of the data (1), Assess the data with a model (2), Model and assess the same data (3). |
| <b>Training Fraction</b><br>( <i>TrainingFraction</i> )       | Specify the fraction of values from the input dataset to be used for model fitting. The exact set of values is chosen at random from the dataset.   | 0.1 | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Probe Location

Sample data attributes at the points in a point cloud.

The Probe filter samples the data set attributes of the current data set at the points in a point cloud. The Probe filter uses interpolation to determine the values at the selected point, whether or not it lies at an input point. The Probe filter operates on any type of data and produces polygonal output (a point cloud).

| Property                               | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )       | This property specifies the dataset from which to obtain probe values.                             |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point or cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet, vtkCompositeDataSet. |
| <b>Probe Type</b><br>( <i>Source</i> ) | This property specifies the dataset whose geometry will be used in determining positions to probe. |                  | The selected object must be the result of the following: sources (includes readers).<br><br>The value must be set to one of the following:<br>FixedRadiusPointSource.  |

## Process Id Scalars

This filter uses colors to show how data is partitioned across processes.

The Process Id Scalars filter assigns a unique scalar value to each piece of the input according to which processor it resides on. This filter operates on any type of data when ParaView is run in parallel. It is useful for determining whether your data is load-balanced across the processors being used. The output data set type is the same as that of the input.

| Property                                    | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )            | This property specifies the input to the Process Id Scalars filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Random Mode</b><br>( <i>RandomMode</i> ) | The value of this property determines whether to use random id values for the various pieces. If set to 1, the unique value per piece will be chosen at random; otherwise the unique value will match the id of the process. | 0                | Only the values 0 and 1 are accepted.  |

## Programmable Filter

Executes a user supplied python script on its input dataset to produce an output dataset.

This filter will execute a python script to produce an output dataset.

The filter keeps a copy of the python script in Script, and creates

Interpreter, a python interpreter to run the script upon the first execution.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Copy Arrays</b><br>( <i>CopyArrays</i> )                      | If this property is set to true, all the cell and point arrays from first input are copied to the output.   | 0                | Only the values 0 and 1 are accepted.  |
| <b>RequestInformation Script</b><br>( <i>InformationScript</i> ) | This property is a python script that is executed during the RequestInformation pipeline pass. Use this to provide information such as WHOLE_EXTENT to the pipeline downstream. |                  |  |
| <b>Input</b><br>( <i>Input</i> )                                 | This property specifies the input(s) to the programmable filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.   |
| <b>Output Data Set Type</b><br>( <i>OutputDataSetType</i> )      | The value of this property determines the dataset type for the output of the programmable filter.   | 8                | The value must be one of the following: Same as Input (8), vtkPolyData (0), vtkStructuredGrid (2), vtkRectilinearGrid (3), vtkUnstructuredGrid (4), vtkImageData (6), vtkUniformGrid (10), vtkMultiblockDataSet (13), vtkHierarchicalBoxDataSet (15), vtkTable (19). |
| <b>Python Path</b><br>( <i>PythonPath</i> )                      | A semi-colon (;) separated list of directories to add to the python library search path.  |                  |  |
| <b>Script</b><br>( <i>Script</i> )                               | This property contains the text of a python program that the programmable filter runs.  |                  |  |

|  |  |  |  |
|--|--|--|--|
| <b>RequestUpdateExtent Script</b><br>( <i>UpdateExtentScript</i> ) | This property is a python script that is executed during the RequestUpdateExtent pipeline pass. Use this to modify the update extent that your filter ask up stream for. |  |  |
|--|--|--|--|

## Python Calculator

This filter evaluates a Python expression

This filter uses Python to calculate an expression.

It depends heavily on the numpy and paraview.vtk modules.

To use the parallel functions, mpi4py is also necessary. The expression is evaluated and the resulting scalar value or numpy array is added to the output as an array. See numpy and paraview.vtk documentation for the list of available functions.

This filter tries to make it easy for the user to write expressions by defining certain variables. The filter tries to assign each array to a variable of the same name. If the name of the array is not a valid Python variable, it has to be accessed through a dictionary called arrays (i.e. arrays['array\_name']). The points can be accessed using the points variable.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Array Association</b><br>( <i>ArrayAssociation</i> ) | This property controls the association of the output array as well as which arrays are defined as variables. | 0                | The value must be one of the following: Point Data (0), Cell Data (1).   |
| <b>Array Name</b><br>( <i>ArrayName</i> )               | The name of the output array.  | result           |  |
| <b>Copy Arrays</b><br>( <i>CopyArrays</i> )             | If this property is set to true, all the cell and point arrays from first input are copied to the output.    | 1                | Only the values 0 and 1 are accepted.  |
| <b>Expression</b><br>( <i>Expression</i> )              | The Python expression evaluated during execution.  |                  |  |
| <b>Input</b><br>( <i>Input</i> )                        | Set the input of the filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Quadric Clustering

This filter is the same filter used to generate level of detail for ParaView. It uses a structured grid of bins and merges all points contained in each bin.

The Quadric Clustering filter produces a reduced-resolution polygonal approximation of the input polygonal dataset. This filter is the one used by ParaView for computing LODs. It uses spatial binning to reduce the number of points in the data set; points that lie within the same spatial bin are collapsed into one representative point.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Copy Cell Data</b><br>( <i>CopyCellData</i> )                 | If this property is set to 1, the cell data from the input will be copied to the output.  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                                 | This property specifies the input to the Quadric Clustering filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Number of Dimensions</b><br>( <i>NumberOfDivisions</i> )      | This property specifies the number of bins along the X, Y, and Z axes of the data set.  | 50 50 50         |   |
| <b>Use Feature Edges</b><br>( <i>UseFeatureEdges</i> )           | If this property is set to 1, feature edge quadrics will be used to maintain the boundary edges along processor divisions.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Use Feature Points</b><br>( <i>UseFeaturePoints</i> )         | If this property is set to 1, feature point quadrics will be used to maintain the boundary points along processor divisions.  | 0                | Only the values 0 and 1 are accepted.   |
| <b>Use Input Points</b><br>( <i>UseInputPoints</i> )             | If the value of this property is set to 1, the representative point for each bin is selected from one of the input points that lies in that bin; the input point that produces the least error is chosen. If the value of this property is 0, the location of the representative point is calculated to produce the least error possible for that bin, but the point will most likely not be one of the input points. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Use Internal Triangles</b><br>( <i>UseInternalTriangles</i> ) | If this property is set to 1, triangles completely contained in a spatial bin will be included in the computation of the bin's quadrics. When this property is set to 0, the filters operates faster, but the resulting surface may not be as well-behaved.   | 0                | Only the values 0 and 1 are accepted.   |

## Random Vectors

This filter creates a new 3-component point data array and sets it as the default vector array. It uses a random number generator to create values.

The Random Vectors filter generates a point-centered array of random vectors. It uses a random number generator to determine the components of the vectors. This filter operates on any type of data set, and the output data set will be of the same type as the input.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Random Vectors filter.                   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Maximum Speed</b><br>( <i>MaximumSpeed</i> ) | This property specifies the maximum length of the random point vectors generated. | 1                | The value must be greater than or equal to 0.  |
| <b>Minimum Speed</b><br>( <i>MinimumSpeed</i> ) | This property specifies the minimum length of the random point vectors generated. | 0                | The value must be greater than or equal to 0.  |

## Rectilinear Grid Connectivity

Parallel fragments extraction and attributes integration on rectilinear grids.

Extracts material fragments from multi-block `vtkRectilinearGrid` datasets based on the selected volume fraction array(s) and a fraction isovalue and integrates the associated attributes.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Double Volume Arrays</b><br>( <i>AddDoubleVolumeArrayName</i> )                   | This property specifies the name(s) of the volume fraction array(s) for generating parts. |                  | An array of scalars is required.  |
| <b>Float Volume Arrays</b><br>( <i>AddFloatVolumeArrayName</i> )                     | This property specifies the name(s) of the volume fraction array(s) for generating parts. |                  | An array of scalars is required.  |
| <b>Unsigned Character Volume Arrays</b><br>( <i>AddUnsignedCharVolumeArrayName</i> ) | This property specifies the name(s) of the volume fraction array(s) for generating parts. |                  | An array of scalars is required.  |
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the input of the filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a cell array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkRectilinearGrid</code> , <code>vtkCompositeDataSet</code> . |
| <b>Volume Fraction Value</b><br>( <i>VolumeFractionSurfaceValue</i> )                | The value of this property is the volume fraction value for the surface.                  | 0.1              | The value must be greater than or equal to 0 and less than or equal to 1.   |

## Reflect

This filter takes the union of the input and its reflection over an axis-aligned plane.

The Reflect filter reflects the input dataset across the specified plane. This filter operates on any type of data set and produces an unstructured grid output.

| Property                                  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>Center</b><br>( <i>Center</i> )        | If the value of the Plane property is X, Y, or Z, then the value of this property specifies the center of the reflection plane.   | 0                |  |
| <b>Copy Input</b><br>( <i>CopyInput</i> ) | If this property is set to 1, the output will contain the union of the input dataset and its reflection. Otherwise the output will contain only the reflection of the input data. | 1                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )          | This property specifies the input to the Reflect filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): <code>vtkDataSet</code> . |

|                                  |  |   |  |
|----------------------------------|--|---|--|
| <b>Plane</b><br>( <i>Plane</i> ) | The value of this property determines which plane to reflect across. If the value is X, Y, or Z, the value of the Center property determines where the plane is placed along the specified axis. The other six options (X Min, X Max, etc.) place the reflection plane at the specified face of the bounding box of the input dataset. | 0 | The value must be one of the following: X Min (0), Y Min (1), Z Min (2), X Max (3), Y Max (4), Z Max (5), X (6), Y (7), Z (8). |
|----------------------------------|--|---|--|

## Resample With Dataset

Sample data attributes at the points of a dataset.

Probe is a filter that computes point attributes at specified point positions. The filter has two inputs: the Input and Source. The Input geometric structure is passed through the filter. The point attributes are computed at the Input point positions by interpolating into the source data. For example, we can compute data values on a plane (plane specified as Input) from a volume (Source). The cell data of the source data is copied to the output based on in which source cell each input point is. If an array of the same name exists both in source's point and cell data, only the one from the point data is probed.

| Property                           | Description  | Default Value(s) | Restrictions   |
|------------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )   | This property specifies the dataset from which to obtain probe values.                             |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point or cell array.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet, vtkCompositeDataSet. |
| <b>Source</b><br>( <i>Source</i> ) | This property specifies the dataset whose geometry will be used in determining positions to probe. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet.   |

## Ribbon

This filter generates ribbon surface from lines. It is useful for displaying streamlines.

The Ribbon filter creates ribbons from the lines in the input data set. This filter is useful for visualizing streamlines. Both the input and output of this filter are polygonal data. The input data set must also have at least one point-centered vector array.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Angle</b><br>( <i>Angle</i> )                  | The value of this property specifies the offset angle (in degrees) of the ribbon from the line normal.   | 0                | The value must be greater than or equal to 0 and less than or equal to 360.   |
| <b>Default Normal</b><br>( <i>DefaultNormal</i> ) | The value of this property specifies the normal to use when the UseDefaultNormal property is set to 1 or the input contains no vector array (SelectInputVectors property). | 0 0 1            |   |
| <b>Input</b><br>( <i>Input</i> )                  | This property specifies the input to the Ribbon filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |

|  |  |   |  |
|--|--|---|--|
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )          | The value of this property indicates the name of the input scalar array used by this filter. The width of the ribbons will be varied based on the values in the specified array if the value of the Width property is 1.   |   | An array of scalars is required.   |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )          | The value of this property indicates the name of the input vector array used by this filter. If the UseDefaultNormal property is set to 0, the normal vectors for the ribbons come from the specified vector array.  | 1 | An array of vectors is required.   |
| <b>Use Default Normal</b><br>( <i>UseDefaultNormal</i> ) | If this property is set to 0, and the input contains no vector array, then default ribbon normals will be generated (DefaultNormal property); if a vector array has been set (SelectInputVectors property), the ribbon normals will be set from the specified array. If this property is set to 1, the default normal (DefaultNormal property) will be used, regardless of whether the SelectInputVectors property has been set. | 0 | Only the values 0 and 1 are accepted.  |
| <b>Vary Width</b><br>( <i>VaryWidth</i> )                | If this property is set to 1, the ribbon width will be scaled according to the scalar array specified in the SelectInputScalars property. Toggle the variation of ribbon width with scalar value.  | 0 | Only the values 0 and 1 are accepted.  |
| <b>Width</b><br>( <i>Width</i> )                         | If the VaryWidth property is set to 1, the value of this property is the minimum ribbon width. If the VaryWidth property is set to 0, the value of this property is half the width of the ribbon.  | 1 | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 0.01. |

## Rotational Extrusion

This filter generates a swept surface while translating the input along a circular path.

The Rotational Extrusion filter forms a surface by rotating the input about the Z axis. This filter is intended to operate on 2D polygonal data. It produces polygonal output.

| Property                                      | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Angle</b><br>( <i>Angle</i> )              | This property specifies the angle of rotation in degrees. The surface is swept from 0 to the value of this property.   | 360              |  |
| <b>Capping</b><br>( <i>Capping</i> )          | If this property is set to 1, the open ends of the swept surface will be capped with a copy of the input dataset. This works property if the input is a 2D surface composed of filled polygons. If the input dataset is a closed solid (e.g., a sphere), then either two copies of the dataset will be drawn or no surface will be drawn. No surface is drawn if either this property is set to 0 or if the two surfaces would occupy exactly the same 3D space (i.e., the Angle property's value is a multiple of 360, and the values of the Translation and DeltaRadius properties are 0). | 1                | Only the values 0 and 1 are accepted.  |
| <b>Delta Radius</b><br>( <i>DeltaRadius</i> ) | The value of this property specifies the change in radius during the sweep process.  | 0                |  |
| <b>Input</b><br>( <i>Input</i> )              | This property specifies the input to the Rotational Extrusion filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkPolyData. |
| <b>Resolution</b><br>( <i>Resolution</i> )    | The value of this property controls the number of intermediate node points used in performing the sweep (rotating from 0 degrees to the value specified by the Angle property).  | 12               | The value must be greater than or equal to 1.  |
| <b>Translation</b><br>( <i>Translation</i> )  | The value of this property specifies the total amount of translation along the Z axis during the sweep process. Specifying a non-zero value for this property allows you to create a corkscrew (value of DeltaRadius > 0) or spring effect.  | 0                |  |



## Scatter Plot

Creates a scatter plot from a dataset.

This filter creates a scatter plot from a dataset. In point data mode, it uses the X point coordinates as the default X array. All other arrays are passed to the output and can be used in the scatter plot. In cell data mode, the first single component array is used as the default X array.

| Property                         | Description                                      | Default Value(s) | Restrictions   |
|----------------------------------|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Shrink

This filter shrinks each input cell so they pull away from their neighbors.

The Shrink filter causes the individual cells of a dataset to break apart from each other by moving each cell's points toward the centroid of the cell. (The centroid of a cell is the average position of its points.) This filter operates on any type of dataset and produces unstructured grid output.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Shrink filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Shrink Factor</b><br>( <i>ShrinkFactor</i> ) | The value of this property determines how far the points will move. A value of 0 positions the points at the centroid of the cell; a value of 1 leaves them at their original positions. | 0.5              | The value must be greater than or equal to 0 and less than or equal to 1.  |

## Slice

This filter slices a data set with a plane. Slicing is similar to a contour. It creates surfaces from volumes and lines from surfaces.

This filter extracts the portion of the input dataset that lies along the specified plane. The Slice filter takes any type of dataset as input. The output of this filter is polygonal data.

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Slice Offset Values</b><br>( <i>ContourValues</i> ) | The values in this property specify a list of current offset values. This can be used to create multiple slices with different centers. Each entry represents a new slice with its center shifted by the offset value. |                  | Determine the length of the dataset's diagonal. The value must lie within -diagonal length to +diagonal length.  |
| <b>Slice Type</b><br>( <i>CutFunction</i> )            | This property sets the parameters of the slice function.   |                  | The value must be set to one of the following: Plane, Box, Sphere.   |
| <b>Input</b><br>( <i>Input</i> )                       | This property specifies the input to the Slice filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Smooth

This filter smooths a polygonal surface by iteratively moving points toward their neighbors.

The Smooth filter operates on a polygonal data set by iteratively adjusting the position of the points using Laplacian smoothing. (Because this filter only adjusts point positions, the output data set is also polygonal.) This results in better-shaped cells and more evenly distributed points.

The Convergence slider limits the maximum motion of any point. It is expressed as a fraction of the length of the diagonal of the bounding box of the data set. If the maximum point motion during a smoothing iteration is less than the Convergence value, the smoothing operation terminates.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Convergence</b><br>( <i>Convergence</i> )                 | The value of this property limits the maximum motion of any point. It is expressed as a fraction of the length of the diagonal of the bounding box of the input dataset. If the maximum point motion during a smoothing iteration is less than the value of this property, the smoothing operation terminates. | 0                | The value must be greater than or equal to 0 and less than or equal to 1.   |
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Smooth filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Number of Iterations</b><br>( <i>NumberOfIterations</i> ) | This property sets the maximum number of smoothing iterations to perform. More iterations produce better smoothing.  | 20               | The value must be greater than or equal to 0.   |

## Stream Tracer

Integrate streamlines in a vector field.

The Stream Tracer filter generates streamlines in a vector field from a collection of seed points. Production of streamlines terminates if a streamline crosses the exterior boundary of the input dataset. Other reasons for termination are listed for the MaximumNumberOfSteps, TerminalSpeed, and MaximumPropagation properties. This filter operates on any type of dataset, provided it has point-centered vectors. The output is polygonal data containing polylines.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Initial Step Length</b><br>( <i>InitialIntegrationStep</i> )   | This property specifies the initial integration step size. For non-adaptive integrators (Runge-Kutta 2 and Runge-Kutta 4), it is fixed (always equal to this initial value) throughout the integration. For an adaptive integrator (Runge-Kutta 4-5), the actual step size varies such that the numerical error is less than a specified threshold. | 0.2              |   |
| <b>Input</b><br>( <i>Input</i> )                                  | This property specifies the input to the Stream Tracer filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 3 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Integration Direction</b><br>( <i>IntegrationDirection</i> )   | This property determines in which direction(s) a streamline is generated.   | 2                | The value must be one of the following: FORWARD (0), BACKWARD (1), BOTH (2).  |
| <b>Integration Step Unit</b><br>( <i>IntegrationStepUnit</i> )    | This property specifies the unit for Minimum/Initial/Maximum integration step size. The Length unit refers to the arc length that a particle travels/advects within a single step. The Cell Length unit represents the step size as a number of cells.  | 2                | The value must be one of the following: Length (1), Cell Length (2).  |
| <b>Integrator Type</b><br>( <i>IntegratorType</i> )               | This property determines which integrator (with increasing accuracy) to use for creating streamlines.   | 2                | The value must be one of the following: Runge-Kutta 2 (0), Runge-Kutta 4 (1), Runge-Kutta 4-5 (2).  |
| <b>Interpolator Type</b><br>( <i>InterpolatorType</i> )           | This property determines which interpolator to use for evaluating the velocity vector field. The first is faster though the second is more robust in locating cells during streamline integration.  | 0                | The value must be one of the following: Interpolator with Point Locator (0), Interpolator with Cell Locator (1).  |
| <b>Maximum Error</b><br>( <i>MaximumError</i> )                   | This property specifies the maximum error (for Runge-Kutta 4-5) tolerated throughout streamline integration. The Runge-Kutta 4-5 integrator tries to adjust the step size such that the estimated error is less than this threshold.  | 1e-06            |   |
| <b>Maximum Step Length</b><br>( <i>MaximumIntegrationStep</i> )   | When using the Runge-Kutta 4-5 ingrator, this property specifies the maximum integration step size.   | 0.5              |   |
| <b>Maximum Steps</b><br>( <i>MaximumNumberOfSteps</i> )           | This property specifies the maximum number of steps, beyond which streamline integration is terminated.   | 2000             |   |
| <b>Maximum Streamline Length</b><br>( <i>MaximumPropagation</i> ) | This property specifies the maximum streamline length (i.e., physical arc length), beyond which line integration is terminated.   | 1                | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 1.   |
| <b>Minimum Step Length</b><br>( <i>MinimumIntegrationStep</i> )   | When using the Runge-Kutta 4-5 ingrator, this property specifies the minimum integration step size.   | 0.01             |   |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )                   | This property contains the name of the vector array from which to generate streamlines.   |                  | An array of vectors is required.  |

|   |   |       |  |
|---|---|-------|--|
| <b>Seed Type</b><br>( <i>Source</i> )             | The value of this property determines how the seeds for the streamlines will be generated.            |       | The selected object must be the result of the following: sources (includes readers).<br><br>The value must be set to one of the following: PointSource, HighResLineSource. |
| <b>Terminal Speed</b><br>( <i>TerminalSpeed</i> ) | This property specifies the terminal speed, below which particle advection/integration is terminated. | 1e-12 |  |

## Stream Tracer With Custom Source

Integrate streamlines in a vector field.

The Stream Tracer filter generates streamlines in a vector field from a collection of seed points. Production of streamlines terminates if a streamline crosses the exterior boundary of the input dataset. Other reasons for termination are listed for the MaximumNumberOfSteps, TerminalSpeed, and MaximumPropagation properties. This filter operates on any type of dataset, provided it has point-centered vectors. The output is polygonal data containing polylines. This filter takes a Source input that provides the seed points.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Initial Step Length</b><br>( <i>InitialIntegrationStep</i> ) | This property specifies the initial integration step size. For non-adaptive integrators (Runge-Kutta 2 and Runge-Kutta 4), it is fixed (always equal to this initial value) throughout the integration. For an adaptive integrator (Runge-Kutta 4-5), the actual step size varies such that the numerical error is less than a specified threshold. | 0.2              |   |
| <b>Input</b><br>( <i>Input</i> )                                | This property specifies the input to the Stream Tracer filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 3 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Integration Direction</b><br>( <i>IntegrationDirection</i> ) | This property determines in which direction(s) a streamline is generated.   | 2                | The value must be one of the following: FORWARD (0), BACKWARD (1), BOTH (2).  |
| <b>Integration Step Unit</b><br>( <i>IntegrationStepUnit</i> )  | This property specifies the unit for Minimum/Initial/Maximum integration step size. The Length unit refers to the arc length that a particle travels/adveects within a single step. The Cell Length unit represents the step size as a number of cells.   | 2                | The value must be one of the following: Length (1), Cell Length (2).  |
| <b>Integrator Type</b><br>( <i>IntegratorType</i> )             | This property determines which integrator (with increasing accuracy) to use for creating streamlines.   | 2                | The value must be one of the following: Runge-Kutta 2 (0), Runge-Kutta 4 (1), Runge-Kutta 4-5 (2).  |
| <b>Maximum Error</b><br>( <i>MaximumError</i> )                 | This property specifies the maximum error (for Runge-Kutta 4-5) tolerated throughout streamline integration. The Runge-Kutta 4-5 integrator tries to adjust the step size such that the estimated error is less than this threshold.  | 1e-06            |   |

|   |   |       |   |
|---|---|-------|---|
| <b>Maximum Step Length</b><br>( <i>MaximumIntegrationStep</i> )   | When using the Runge-Kutta 4-5 ingrator, this property specifies the maximum integration step size.                             | 0.5   |   |
| <b>Maximum Steps</b><br>( <i>MaximumNumberOfSteps</i> )           | This property specifies the maximum number of steps, beyond which streamline integration is terminated.                         | 2000  |   |
| <b>Maximum Streamline Length</b><br>( <i>MaximumPropagation</i> ) | This property specifies the maximum streamline length (i.e., physical arc length), beyond which line integration is terminated. | 1     | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 1. |
| <b>Minimum Step Length</b><br>( <i>MinimumIntegrationStep</i> )   | When using the Runge-Kutta 4-5 ingrator, this property specifies the minimum integration step size.                             | 0.01  |   |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )                   | This property contains the name of the vector array from which to generate streamlines.   |       | An array of vectors is required.  |
| <b>Source</b><br>( <i>Source</i> )                                | This property specifies the input used to obtain the seed points.   |       | The selected object must be the result of the following: sources (includes readers).                |
| <b>Terminal Speed</b><br>( <i>TerminalSpeed</i> )                 | This property specifies the terminal speed, below which particle advection/integration is terminated.                           | 1e-12 |   |

## Subdivide

This filter iteratively divide triangles into four smaller triangles. New points are placed linearly so the output surface matches the input surface.

The Subdivide filter iteratively divides each triangle in the input dataset into 4 new triangles. Three new points are added per triangle -- one at the midpoint of each edge. This filter operates only on polygonal data containing triangles, so run your polygonal data through the Triangulate filter first if it is not composed of triangles. The output of this filter is also polygonal.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                                 | This parameter specifies the input to the Subdivide filter.                           |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Number of Subdivisions</b><br>( <i>NumberOfSubdivisions</i> ) | The value of this property specifies the number of subdivision iterations to perform. | 1                | The value must be greater than or equal to 1 and less than or equal to 4.   |

## Surface Flow

This filter integrates flow through a surface.

The flow integration filter integrates the dot product of a point flow vector field and surface normal. It computes the net flow across the 2D surface. It operates on any type of dataset and produces an unstructured grid output.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Surface Flow filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 3 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Select Input Vectors</b><br>( <i>SelectInputVectors</i> ) | The value of this property specifies the name of the input vector array containing the flow vector field. |                  | An array of vectors is required.  |

## Surface Vectors

This filter constrains vectors to lie on a surface.

The Surface Vectors filter is used for 2D data sets. It constrains vectors to lie in a surface by removing components of the vectors normal to the local surface.

| Property   | Description   | Default Value(s) | Restrictions  |
|--|---|------------------|---|
| <b>Constraint Mode</b><br>( <i>ConstraintMode</i> )          | This property specifies whether the vectors will be parallel or perpendicular to the surface. If the value is set to PerpendicularScale (2), then the output will contain a scalar array with the dot product of the surface normal and the vector at each point. | 0                | The value must be one of the following: Parallel (0), Perpendicular (1), PerpendicularScale (2).  |
| <b>Input</b><br>( <i>Input</i> )                             | This property specifies the input to the Surface Vectors filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 3 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Select Input Vectors</b><br>( <i>SelectInputVectors</i> ) | This property specifies the name of the input vector array to process.  |                  | An array of vectors is required.  |

## Table To Points

Converts table to set of points.

The TableToPolyData filter converts a vtkTable to a set of points in a vtkPolyData. One must specify the columns in the input table to use as the X, Y and Z coordinates for the points in the output.

| Property                              | Description   | Default Value(s) | Restrictions  |
|---------------------------------------|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )      | This property specifies the input..   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable. |
| <b>X Column</b><br>( <i>XColumn</i> ) | This property specifies which data array is going to be used as the X coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |
| <b>Y Column</b><br>( <i>YColumn</i> ) | This property specifies which data array is going to be used as the Y coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |
| <b>Z Column</b><br>( <i>ZColumn</i> ) | This property specifies which data array is going to be used as the Z coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |

## Table To Structured Grid

Converts to table to structured grid.

The TableToStructuredGrid filter converts a vtkTable to a vtkStructuredGrid. One must specifies the columns in the input table to use as the X, Y and Z coordinates for the points in the output, and the whole extent.

| Property                                      | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )              | This property specifies the input..   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable. |
| <b>Whole Extent</b><br>( <i>WholeExtent</i> ) |   | 0 0 0 0 0 0      |   |
| <b>X Column</b><br>( <i>XColumn</i> )         | This property specifies which data array is going to be used as the X coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |
| <b>Y Column</b><br>( <i>YColumn</i> )         | This property specifies which data array is going to be used as the Y coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |
| <b>Z Column</b><br>( <i>ZColumn</i> )         | This property specifies which data array is going to be used as the Z coordinate in the generated polydata dataset. |                  | An array of scalars is required.  |

## Temporal Cache

Saves a copy of the data set for a fixed number of time steps.

The Temporal Cache can be used to save multiple copies of a data set at different time steps to prevent thrashing in the pipeline caused by downstream filters that adjust the requested time step. For example, assume that there is a downstream Temporal Interpolator filter. This filter will (usually) request two time steps from the upstream filters, which in turn (usually) causes the upstream filters to run twice, once for each time step. The next time the interpolator requests the same two time steps, they might force the upstream filters to re-evaluate the same two time steps. The Temporal Cache can keep copies of both of these time steps and provide the requested data without

having to run upstream filters.

| Property                                  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Cache Size</b><br>( <i>CacheSize</i> ) | The cache size determines the number of time steps that can be cached at one time. The maximum number is 10. The minimum is 2 (since it makes little sense to cache less than that). | 2                | The value must be greater than or equal to 2 and less than or equal to 10.  |
| <b>Input</b><br>( <i>Input</i> )          | This property specifies the input of the Temporal Cache filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |

## Temporal Interpolator

Interpolate between time steps.

The Temporal Interpolator converts data that is defined at discrete time steps to one that is defined over a continuum of time by linearly interpolating the data's field data between two adjacent time steps. The interpolated values are a simple approximation and should not be interpreted as anything more. The Temporal Interpolator assumes that the topology between adjacent time steps does not change.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Discrete Time Step Interval</b><br>( <i>DiscreteTimeStepInterval</i> ) | If Discrete Time Step Interval is set to 0, then the Temporal Interpolator will provide a continuous region of time on its output. If set to anything else, then the output will define a finite set of time points on its output, each spaced by the Discrete Time Step Interval. The output will have (time range)/(discrete time step interval) time steps. (Note that the time range is defined by the time range of the data of the input filter, which may be different from other pipeline objects or the range defined in the animation inspector.) This is a useful option to use if you have a dataset with one missing time step and wish to 'file-in' the missing data with an interpolated value from the steps on either side. | 0                |   |
| <b>Input</b><br>( <i>Input</i> )  | This property specifies the input of the Temporal Interpolator.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |



## Temporal Shift Scale

Shift and scale time values.

The Temporal Shift Scale filter linearly transforms the time values of a pipeline object by applying a shift and then scale. Given a data at time  $t$  on the input, it will be transformed to time  $t \cdot \text{Shift} + \text{Scale}$  on the output. Inversely, if this filter has a request for time  $t$ , it will request time  $(t - \text{Shift}) / \text{Scale}$  on its input.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                                      | The input to the Temporal Shift Scale filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |
| <b>Maximum Number Of Periods</b><br>( <i>MaximumNumberOfPeriods</i> ) | If Periodic time is enabled, this controls how many time periods time is reported for. A filter cannot output an infinite number of time steps and therefore a finite number of periods is generated when reporting time.  | 1                | The value must be greater than or equal to 0 and less than or equal to 100.   |
| <b>Periodic</b><br>( <i>Periodic</i> )                                | If Periodic is true, requests for time will be wrapped around so that the source appears to be a periodic time source. If data exists for times $\{0, N-1\}$ , setting periodic to true will cause time 0 to be produced when time $N$ , $2N$ , $2N$ etc is requested. This effectively gives the source the ability to generate time data indefinitely in a loop. When combined with Shift/Scale, the time becomes periodic in the shifted and scaled time frame of reference. Note: Since the input time may not start at zero, the wrapping of time from the end of one period to the start of the next, will subtract the initial time - a source with $T\{5..6\}$ repeated periodically will have output time $\{5..6..7..8\}$ etc. | 0                | Only the values 0 and 1 are accepted.   |
| <b>Periodic End Correction</b><br>( <i>PeriodicEndCorrection</i> )    | If Periodic time is enabled, this flag determines if the last time step is the same as the first. If PeriodicEndCorrection is true, then it is assumed that the input data goes from 0-1 (or whatever scaled/shifted actual time) and time 1 is the same as time 0 so that steps will be 0,1,2,3... $N$ ,1,2,3... $N$ ,1,2,3 where step $N$ is the same as 0 and step 0 is not repeated. When this flag is false the data is assumed to be literal and output is of the form 0,1,2,3... $N$ ,0,1,2,3... By default this flag is ON   | 1                | Only the values 0 and 1 are accepted.   |
| <b>Post Shift</b><br>( <i>PostShift</i> )                             | The amount of time the input is shifted.   | 0                |   |
| <b>Pre Shift</b><br>( <i>PreShift</i> )                               | Apply a translation to the data before scaling. To convert $T\{5,100\}$ to $T\{0,1\}$ use $\text{Preshift}=-5$ , $\text{Scale}=1/95$ , $\text{PostShift}=0$ To convert $T\{5,105\}$ to $T\{5,10\}$ use $\text{Preshift}=-5$ , $\text{Scale}=5/100$ , $\text{PostShift}=5$  | 0                |   |
| <b>Scale</b><br>( <i>Scale</i> )                                      | The factor by which the input time is scaled.  | 1                |   |

## Temporal Snap-to-Time-Step

Modifies the time range/steps of temporal data.

This filter modifies the time range or time steps of the data without changing the data itself. The data is not resampled by this filter, only the information accompanying the data is modified.

| Property                                | Description                                      | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )        | This property specifies the input of the filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |
| <b>Snap Mode</b><br>( <i>SnapMode</i> ) | Determine which time step to snap to.            | 0                | The value must be one of the following: Nearest (0), NextBelowOrEqual (1), NextAboveOrEqual (2).  |

## Temporal Statistics

Loads in all time steps of a data set and computes some statistics about how each point and cell variable changes over time.

Given an input that changes over time, vtkTemporalStatistics looks at the data for each time step and computes some statistical information of how a point or cell variable changes over time. For example, vtkTemporalStatistics can compute the average value of "pressure" over time of each point.

Note that this filter will require the upstream filter to be run on every time step that it reports that it can compute. This may be a time consuming operation.

vtkTemporalStatistics ignores the temporal spacing. Each timestep will be weighted the same regardless of how long of an interval it is to the next timestep. Thus, the average statistic may be quite different from an integration of the variable if the time spacing varies.

| Property   | Description   | Default Value(s) | Restrictions   |
|--|---|------------------|--|
| <b>Compute Average</b><br>( <i>ComputeAverage</i> )                      | Compute the average of each point and cell variable over time.            | 1                | Only the values 0 and 1 are accepted.  |
| <b>Compute Maximum</b><br>( <i>ComputeMaximum</i> )                      | Compute the maximum of each point and cell variable over time.            | 1                | Only the values 0 and 1 are accepted.  |
| <b>Compute Minimum</b><br>( <i>ComputeMinimum</i> )                      | Compute the minimum of each point and cell variable over time.            | 1                | Only the values 0 and 1 are accepted.  |
| <b>Compute Standard Deviation</b><br>( <i>ComputeStandardDeviation</i> ) | Compute the standard deviation of each point and cell variable over time. | 1                | Only the values 0 and 1 are accepted.  |
| <b>Input</b><br>( <i>Input</i> )   | Set the input to the Temporal Statistics filter.                          |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Tessellate

Tessellate nonlinear curves, surfaces, and volumes with lines, triangles, and tetrahedra.

The Tessellate filter tessellates cells with nonlinear geometry and/or scalar fields into a simplicial complex with linearly interpolated field values that more closely approximate the original field. This is useful for datasets containing quadratic cells.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>Chord Error</b><br>( <i>ChordError</i> )                                     | This property controls the maximum chord error allowed at any edge midpoint in the output tessellation. The chord error is measured as the distance between the midpoint of any output edge and the original nonlinear geometry.   | 0.001            |   |
| <b>Field Error</b><br>( <i>FieldError2</i> )                                    | This property controls the maximum field error allowed at any edge midpoint in the output tessellation. The field error is measured as the difference between a field value at the midpoint of an output edge and the value of the corresponding field in the original nonlinear geometry.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | This property specifies the input to the Tessellate filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkPolyData, vtkDataSet, vtkUnstructuredGrid. |
| <b>Maximum Number of Subdivisions</b><br>( <i>MaximumNumberOfSubdivisions</i> ) | This property specifies the maximum number of times an edge may be subdivided. Increasing this number allows further refinement but can drastically increase the computational and storage requirements, especially when the value of the OutputDimension property is 3.   | 3                | The value must be greater than or equal to 0 and less than or equal to 8.   |
| <b>Merge Points</b><br>( <i>MergePoints</i> )                                   | If the value of this property is set to 1, coincident vertices will be merged after tessellation has occurred. Only geometry is considered during the merge and the first vertex encountered is the one whose point attributes will be used. Any discontinuities in point fields will be lost. On the other hand, many operations, such as streamline generation, require coincident vertices to be merged. Toggle whether to merge coincident vertices. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Output Dimension</b><br>( <i>OutputDimension</i> )                           | The value of this property sets the maximum dimensionality of the output tessellation. When the value of this property is 3, 3D cells produce tetrahedra, 2D cells produce triangles, and 1D cells produce line segments. When the value is 2, 3D cells will have their boundaries tessellated with triangles. When the value is 1, all cells except points produce line segments.   | 3                | The value must be greater than or equal to 1 and less than or equal to 3.   |

## Tetrahedralize

This filter converts 3-d cells to tetrahedrons and polygons to triangles. The output is always of type unstructured grid.

The Tetrahedralize filter converts the 3D cells of any type of dataset to tetrahedrons and the 2D ones to triangles. This filter always produces unstructured grid output.

| Property                         | Description   | Default Value(s) | Restrictions   |
|----------------------------------|---|------------------|--|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Tetrahedralize filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## Texture Map to Cylinder

Generate texture coordinates by mapping points to cylinder.

This is a filter that generates 2D texture coordinates by mapping input dataset points onto a cylinder. The cylinder is generated automatically.

The cylinder is generated automatically by computing the axis of the cylinder. Note that the generated texture coordinates for the s-coordinate ranges from (0-1) (corresponding to angle of 0->360 around axis), while the mapping of the t-coordinate is controlled by the projection of points along the axis.

| Property                                      | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )              | Set the input to the Texture Map to Cylinder filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Prevent Seam</b><br>( <i>PreventSeam</i> ) | Control how the texture coordinates are generated. If Prevent Seam is set, the s-coordinate ranges from 0->1 and 1->0 corresponding to the theta angle variation between 0->180 and 180->0 degrees. Otherwise, the s-coordinate ranges from 0->1 between 0->360 degrees. | 1                | Only the values 0 and 1 are accepted.  |

## Texture Map to Plane

Generate texture coordinates by mapping points to plane.

TextureMapToPlane is a filter that generates 2D texture coordinates by mapping input dataset points onto a plane. The plane is generated automatically. A least squares method is used to generate the plane automatically.

| Property | Description | Default Value(s) | Restrictions |
|----------|-------------|------------------|--------------|
|----------|-------------|------------------|--------------|

|                         |   |  |  |
|-------------------------|---|--|--|
| <b>Input</b><br>(Input) | Set the input to the Texture Map to Plane filter. |  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
|-------------------------|---|--|--|

## Texture Map to Sphere

Generate texture coordinates by mapping points to sphere.

This is a filter that generates 2D texture coordinates by mapping input dataset points onto a sphere. The sphere is generated automatically. The sphere is generated automatically by computing the center i.e. averaged coordinates, of the sphere. Note that the generated texture coordinates range between (0,1). The s-coordinate lies in the angular direction around the z-axis, measured counter-clockwise from the x-axis. The t-coordinate lies in the angular direction measured down from the north pole towards the south pole.

| Property                             | Description  | Default Value(s) | Restrictions   |
|--------------------------------------|--|------------------|--|
| <b>Input</b><br>(Input)              | Set the input to the Texture Map to Sphere filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Prevent Seam</b><br>(PreventSeam) | Control how the texture coordinates are generated. If Prevent Seam is set, the s-coordinate ranges from 0->1 and 1->0 corresponding to the theta angle variation between 0->180 and 180->0 degrees. Otherwise, the s-coordinate ranges from 0->1 between 0->360 degrees. | 1                | Only the values 0 and 1 are accepted.  |

## Threshold

This filter extracts cells that have point or cell scalars in the specified range.

The Threshold filter extracts the portions of the input dataset whose scalars lie within the specified range. This filter operates on either point-centered or cell-centered data. This filter operates on any type of dataset and produces unstructured grid output.

To select between these two options, select either Point Data or Cell Data from the Attribute Mode menu. Once the Attribute Mode has been selected, choose the scalar array from which to threshold the data from the Scalars menu. The Lower Threshold and Upper Threshold sliders determine the range of the scalars to retain in the output. The All Scalars check box only takes effect when the Attribute Mode is set to Point Data. If the All Scalars option is checked, then a cell will only be passed to the output if the scalar values of all of its points lie within the range indicated by the Lower Threshold and Upper Threshold sliders. If unchecked, then a cell will be added to the output if the specified scalar value for any of its points is within the chosen range.

| Property  | Description  | Default Value(s) | Restrictions  |
|---|--|------------------|---|
| <b>All Scalars</b><br>( <i>AllScalars</i> )           | If the value of this property is 1, then a cell is only included in the output if the value of the selected array for all its points is within the threshold. This is only relevant when thresholding by a point-centered array. | 1                | Only the values 0 and 1 are accepted.   |
| <b>Input</b><br>( <i>Input</i> )                      | This property specifies the input to the Threshold filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point or cell array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )       | The value of this property contains the name of the scalar array from which to perform thresholding.   |                  | An array of scalars is required.<br><br>Valid array names will be chosen from point and cell data.  |
| <b>Threshold Range</b><br>( <i>ThresholdBetween</i> ) | The values of this property specify the upper and lower bounds of the thresholding operation.  | 0 0              | The value must lie within the range of the selected data array.   |

## Transform

This filter applies transformation to the polygons.

The Transform filter allows you to specify the position, size, and orientation of polygonal, unstructured grid, and curvilinear data sets.

| Property                                 | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> )         | This property specifies the input to the Transform filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |
| <b>Transform</b><br>( <i>Transform</i> ) | The values in this property allow you to specify the transform (translation, rotation, and scaling) to apply to the input dataset. |                  | The selected object must be the result of the following: transforms.<br><br>The value must be set to one of the following: Transform3.  |

## Triangle Strips

This filter uses a greedy algorithm to convert triangles into triangle strips

The Triangle Strips filter converts triangles into triangle strips and lines into polylines. This filter operates on polygonal data sets and produces polygonal output.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>Input</b><br>( <i>Input</i> )                  | This property specifies the input to the Triangle Strips filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Maximum Length</b><br>( <i>MaximumLength</i> ) | This property specifies the maximum number of triangles/lines to include in a triangle strip or polyline. | 1000             | The value must be greater than or equal to 4 and less than or equal to 100000.  |

## Triangulate

This filter converts polygons and triangle strips to basic triangles.

The Triangulate filter decomposes polygonal data into only triangles, points, and lines. It separates triangle strips and polylines into individual triangles and lines, respectively. The output is polygonal data. Some filters that take polygonal data as input require that the data be composed of triangles rather than other polygons, so passing your data through this filter first is useful in such situations. You should use this filter in these cases rather than the Tetrahedralize filter because they produce different output dataset types. The filters referenced require polygonal input, and the Tetrahedralize filter produces unstructured grid output.

| Property                         | Description  | Default Value(s) | Restrictions  |
|----------------------------------|--|------------------|---|
| <b>Input</b><br>( <i>Input</i> ) | This property specifies the input to the Triangulate filter. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |

## Tube

Convert lines into tubes. Normals are used to avoid cracks between tube segments.

The Tube filter creates tubes around the lines in the input polygonal dataset. The output is also polygonal.

| Property   | Description  | Default Value(s) | Restrictions  |
|--|--|------------------|---|
| <b>Capping</b><br>( <i>Capping</i> )               | If this property is set to 1, endcaps will be drawn on the tube. Otherwise the ends of the tube will be open.  | 1                | Only the values 0 and 1 are accepted.   |
| <b>Default Normal</b><br>( <i>DefaultNormal</i> )  | The value of this property specifies the normal to use when the UseDefaultNormal property is set to 1 or the input contains no vector array (SelectInputVectors property). | 0 0 1            |   |
| <b>Input</b><br>( <i>Input</i> )                   | This property specifies the input to the Tube filter.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Number of Sides</b><br>( <i>NumberOfSides</i> ) | The value of this property indicates the number of faces around the circumference of the tube.   | 6                | The value must be greater than or equal to 3.   |

|  |  |    |  |
|--|--|----|--|
| <b>Radius</b><br>( <i>Radius</i> )                       | The value of this property sets the radius of the tube. If the radius is varying (VaryRadius property), then this value is the minimum radius.   | 1  | The value must be less than the largest dimension of the dataset multiplied by a scale factor of 0.01. |
| <b>Radius Factor</b><br>( <i>RadiusFactor</i> )          | If varying the radius (VaryRadius property), the property sets the maximum tube radius in terms of a multiple of the minimum radius. If not varying the radius, this value has no effect.  | 10 |  |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> )          | This property indicates the name of the scalar array on which to operate. The indicated array may be used for scaling the tubes. (See the VaryRadius property.)  |    | An array of scalars is required.   |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> )          | This property indicates the name of the vector array on which to operate. The indicated array may be used for scaling and/or orienting the tubes. (See the VaryRadius property.)   | 1  | An array of vectors is required.   |
| <b>Use Default Normal</b><br>( <i>UseDefaultNormal</i> ) | If this property is set to 0, and the input contains no vector array, then default ribbon normals will be generated (DefaultNormal property); if a vector array has been set (SelectInputVectors property), the ribbon normals will be set from the specified array. If this property is set to 1, the default normal (DefaultNormal property) will be used, regardless of whether the SelectInputVectors property has been set. | 0  | Only the values 0 and 1 are accepted.  |
| <b>Vary Radius</b><br>( <i>VaryRadius</i> )              | The property determines whether/how to vary the radius of the tube. If varying by scalar (1), the tube radius is based on the point-based scalar values in the dataset. If it is varied by vector, the vector magnitude is used in varying the radius.   | 0  | The value must be one of the following: Off (0), By Scalar (1), By Vector (2), By Absolute Scalar (3). |

## Warp By Scalar

This filter moves point coordinates along a vector scaled by a point attribute. It can be used to produce carpet plots.

The Warp (scalar) filter translates the points of the input data set along a vector by a distance determined by the specified scalars. This filter operates on polygonal, curvilinear, and unstructured grid data sets containing single-component scalar arrays. Because it only changes the positions of the points, the output data set type is the same as that of the input. Any scalars in the input dataset are copied to the output, so the data can be colored by them.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Warp (scalar) filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 1 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |
| <b>Normal</b><br>( <i>Normal</i> )              | The values of this property specify the direction along which to warp the dataset if any normals contained in the input dataset are not being used for this purpose. (See the UseNormal property.) | 0 0 1            |  |
| <b>Scale Factor</b><br>( <i>ScaleFactor</i> )   | The scalar value at a given point is multiplied by the value of this property to determine the magnitude of the change vector for that point.  | 1                |  |
| <b>Scalars</b><br>( <i>SelectInputScalars</i> ) | This property contains the name of the scalar array by which to warp the dataset.  |                  | An array of scalars is required.   |



|   |   |   |                                       |
|---|---|---|---------------------------------------|
| <b>Use Normal</b><br>( <i>UseNormal</i> ) | If point normals are present in the dataset, the value of this property toggles whether to use a single normal value (value = 1) or the normals from the dataset (value = 0).                         | 0 | Only the values 0 and 1 are accepted. |
| <b>XY Plane</b><br>( <i>XYPlane</i> )     | If the value of this property is 1, then the Z-coordinates from the input are considered to be the scalar values, and the displacement is along the Z axis. This is useful for creating carpet plots. | 0 | Only the values 0 and 1 are accepted. |

## Warp By Vector

This filter displaces point coordinates along a vector attribute. It is useful for showing mechanical deformation.

The Warp (vector) filter translates the points of the input dataset using a specified vector array. The vector array chosen specifies a vector per point in the input. Each point is translated along its vector by a given scale factor. This filter operates on polygonal, curvilinear, and unstructured grid datasets. Because this filter only changes the positions of the points, the output dataset type is the same as that of the input.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>Input</b><br>( <i>Input</i> )                | This property specifies the input to the Warp (vector) filter.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The dataset must contain a point array with 3 components.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPointSet. |
| <b>Scale Factor</b><br>( <i>ScaleFactor</i> )   | Each component of the selected vector array will be multiplied by the value of this property before being used to compute new point coordinates. | 1                |  |
| <b>Vectors</b><br>( <i>SelectInputVectors</i> ) | The value of this property contains the name of the vector array by which to warp the dataset's point coordinates.                               |                  | An array of vectors is required.   |

# List of Writers

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## AnimationSceneImageWriter

Internal writer to used paraview uses when it disconnects the GUI.

| Property   | Description  | Default Value(s) | Restrictions                                  |
|--|--|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )          | The name of the file to save the animation into.   |                  |   |
| <b>Frame Rate</b><br>( <i>FrameRate</i> )        | Get/Set the frame rate to use for saving the animation. This frame rate is the frame rate that gets saved in the movie file generated, if applicable. If does not affect the FrameRate set on the animation scene at all. In other words, this is the playback frame rate and not the animation generation frame rate. | 0                |   |
| <b>Magnification</b><br>( <i>Magnification</i> ) | The magnification factor to use for the saved animation.   | 1                | The value must be greater than or equal to 1. |

## CSVWriter

Writer to write CSV files from table.

In parallel, it delivers the table to the root node and then saves the CSV. For composite datasets, it saves multiple csv files.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                     | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )                            | The input filter/source whose output dataset is to written to the file.                                       |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkTable. |
| <b>Write All Time Steps</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader. | 0                | Only the values 0 and 1 are accepted.  |

## DataSetCSVWriter

Writer to write CSV files from any dataset. Set FieldAssociation to choose whether cell data/point data needs to be saved.

In parallel, it delivers the table to the root node and then saves the CSV. For composite datasets, it saves multiple csv files.

---

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                     | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )                            | The input filter/source whose output dataset is to be written to the file.                                    |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |
| <b>Write All Time Steps</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader. | 0                | Only the values 0 and 1 are accepted.  |
| <b>Field Association</b><br>( <i>FieldAssociation</i> )     | Select the attribute data to pass.  | 0                | The value must be one of the following: Points (0), Cells (1), Field Data (2).   |

## DataSetWriter

Writer to write any type of data object in a legacy vtk data file.

Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to be written to the file.                                    |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader. | 0                | Only the values 0 and 1 are accepted.   |
| <b>File Type</b><br>( <i>FileType</i> )                                   | The type for the data file (i.e. ascii or binary).  | 2                | The value must be one of the following: Ascii (1), Binary (2).  |

## EnSightWriter

Writer to write unstructured grid data as an EnSight file. Binary files written on one system may not be readable on other systems. Be sure to specify the endian-ness of the file when reading it into EnSight.

| Property                                    | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )     | The name of the file to be written.  |                  |  |
| <b>Ghost Level</b><br>( <i>GhostLevel</i> ) | The number of ghost levels to include in the output file(s). Note that some writers (such as ExodusIIWriter) may not support writing ghost levels. | 0                | The value must be greater than or equal to 0.  |
| <b>Input</b><br>( <i>Input</i> )            | The input filter/source whose output dataset is to be written to the file.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkUnstructuredGrid. |

## ExodusIIWriter

Writer to write Exodus II files. Refere to <http://endo.sandia.gov/SEACAS/> for more information about the Exodus II format.

| Property  | Description  | Default Value(s) | Restrictions   |
|---|--|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                     | The name of the file to be written.  |                  |  |
| <b>Ghost Level</b><br>( <i>GhostLevel</i> )                 | The number of ghost levels to include in the output file(s). Note that some writers (such as ExodusIIWriter) may not support writing ghost levels. | 0                | The value must be greater than or equal to 0.  |
| <b>Input</b><br>( <i>Input</i> )                            | The input filter/source whose output dataset is to be written to the file.   |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them):<br>vtkCompositeDataSet, vtkDataSet. |
| <b>Write All Time Steps</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                                      | 0                | Only the values 0 and 1 are accepted.  |

## MetaImageWriter

Writer to write a binary UNC meta image data. This is a fairly simple yet powerful format consisting of a text header and a binary data section.

MetaImage headers are expected to have extension: ".mha" or ".mhd"

| Property                                | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to be written.                                     |                  |  |
| <b>Input</b><br>( <i>Input</i> )        | The input filter/source whose output dataset is to written to the file. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |

## PDataSetWriter

Writer to write any type of data object in a legacy vtk data file.

This version is used when running in parallel. It gathers data to first node and saves 1 file.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                     | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )                            | The input filter/source whose output dataset is to written to the file.                                       |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataObject. |
| <b>Write All Time Steps</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader. | 0                | Only the values 0 and 1 are accepted.   |
| <b>File Type</b><br>( <i>FileType</i> )                     | The type for the data file (i.e. ascii or binary).  | 2                | The value must be one of the following: Ascii (1), Binary (2).  |

## PNGWriter

Writer to write image data as a PNG file. It supports 1 to 4 component data of unsigned char or unsigned short.

| Property                                | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to be written.                                     |                  |  |
| <b>Input</b><br>( <i>Input</i> )        | The input filter/source whose output dataset is to written to the file. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |

## PPLYWriter

Writer to write polygonal data in Stanford University PLY format. The data can be written in either binary (little or big endian) or ASCII representation. As for PointData and CellData, vtkPLYWriter cannot handle normals or vectors. It only handles RGB PointData and CellData. This version is used when running in parallel. It gathers data to first node and saves 1 file.

| Property                                | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to be written.                                     |                  |   |
| <b>Input</b><br>( <i>Input</i> )        | The input filter/source whose output dataset is to written to the file. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>File Type</b><br>( <i>FileType</i> ) | The type for the data file (i.e. ascii or binary).                      | 2                | The value must be one of the following: Ascii (1), Binary (2).  |

## PSTLWriter

STLWriter writes stereo lithography (.stl) files in either ASCII or binary form. Stereo lithography files only contain triangles. If polygons with more than 3 vertices are present, only the first 3 vertices are written. Use TriangleFilter to convert polygons to triangles. This version of the reader is used when running in parallel. It gathers all the geometry to first node and saves 1 file.

| Property                                | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to be written.                                     |                  |   |
| <b>Input</b><br>( <i>Input</i> )        | The input filter/source whose output dataset is to written to the file. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>File Type</b><br>( <i>FileType</i> ) | The type for the data file (i.e. ascii or binary).                      | 2                | The value must be one of the following: Ascii (1), Binary (2).  |

## XMLDataSetWriterBase

Base for all XML-based file-series writers.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).                  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2). |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLHierarchicalBoxDataWriter

Writer to write a hierarchical box in a xml-based vtk data file. Can be used for parallel writing as well as serial writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkHierarchicalBoxDataSet. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                       |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLHyperOctreeWriter

Writer to write unstructured grid in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to be written to the file.  |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkHyperOctree. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                            |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |

## XMLImageDataWriter

Writer to write image data in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to be written to the file.  |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                          |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |



## XMLMultiBlockDataWriter

Writer to write a multiblock dataset in a xml-based vtk data file. Can be used for parallel writing as well as serial writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkMultiBlockDataSet. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                  |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |

## XMLImageDataWriter

Writer to write image data in a xml-based vtk data file. Can be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkImageData. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                          |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |

## XMLPPolyDataWriter

Writer to write polydata in a xml-based vtk data file. Can be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                         |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLPRectilinearGridWriter

Writer to write rectilinear grid in a xml-based vtk data file. Can be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkRectilinearGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |

## XMLPStructuredGridWriter

Writer to write structured grid in a xml-based vtk data file. Can be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkStructuredGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                               |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLPUnstructuredGridWriter

Writer to write unstructured grid in a xml-based vtk data file. Can be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                 |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLPVAnimationWriter

Internal writer proxy used when saving animation geometry to save all parts of the current source to the file with pieces spread across the server processes.

| Property   | Description  | Default Value(s) | Restrictions |
|--|--|------------------|--------------|
| <b>File Name</b><br>( <i>FileName</i> )              | Name of the file to write.   |                  |              |
| <b>Representations</b><br>( <i>Representations</i> ) | The input filter/source whose output dataset is to be written to the file. |                  |              |

## XMLPVDWriter

Writer to write ParaView data files (pvd). It is used to save all pieces of a source/filter to a file with pieces spread across the server processes.

| Property   | Description  | Default Value(s) | Restrictions   |
|--|--|------------------|--|
| <b>Compressor Type</b><br>( <i>CompressorType</i> )          | The compression algorithm used to compress binary data (appended mode only).   | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                      | The mode used for writing the file's data i.e. ascii, binary, appended binary.   | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).   |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> ) | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected).                | 0                | Only the values 0 and 1 are accepted.  |
| <b>File Name</b><br>( <i>FileName</i> )                      | The name of the file to be written.  |                  |  |
| <b>Ghost Level</b><br>( <i>GhostLevel</i> )                  | The number of ghost levels to include in the output file(s). Note that some writers (such as ExodusIIWriter) may not support writing ghost levels. | 0                | The value must be greater than or equal to 0.  |
| <b>Input</b><br>( <i>Input</i> )                             | The input filter/source whose output dataset is to be written to the files.  |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

## XMLPolyDataWriter

Writer to write poly data in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to be written to the file.  |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkPolyData. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                         |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLRectilinearGridWriter

Writer to write rectilinear grid in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions  |
|---|---|------------------|---|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |   |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to be written to the file.  |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkRectilinearGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.   |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).   |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.   |

## XMLStructuredGridWriter

Writer to write structured grid in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkStructuredGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                               |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XMLUnstructuredGridWriter

Writer to write unstructured grid in a xml-based vtk data file. Cannot be used for parallel writing.

| Property  | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> )                                   | The name of the file to be written.   |                  |  |
| <b>Input</b><br>( <i>Input</i> )  | The input filter/source whose output dataset is to written to the file.   |                  | The selected dataset must be one of the following types (or a subclass of one of them): vtkUnstructuredGrid. |
| <b>Write all timesteps as file-series</b><br>( <i>WriteAllTimeSteps</i> ) | When WriteAllTimeSteps is turned ON, the writer is executed once for each timestep available from the reader.                       | 0                | Only the values 0 and 1 are accepted.  |
| <b>Compressor Type</b><br>( <i>CompressorType</i> )                       | The compression algorithm used to compress binary data (appended mode only).  | 0                | The value must be one of the following: None (0), ZLib (1).  |
| <b>Data Mode</b><br>( <i>DataMode</i> )                                   | The mode uses for writing the file's data i.e. ascii, binary, appended binary.  | 2                | The value must be one of the following: Ascii (0), Binary (1), Appended (2).                                 |
| <b>Encode Appended Data</b><br>( <i>EncodeAppendedData</i> )              | When EncodeAppendedData is turned ON, the writer uses base64 encoding when writing binary data (only if appended mode is selected). | 0                | Only the values 0 and 1 are accepted.  |

## XdmfWriter

Writer to write data in eXtensible Data Model and Format \*(XDMF) files.

| Property                                | Description   | Default Value(s) | Restrictions   |
|---|---|------------------|--|
| <b>File Name</b><br>( <i>FileName</i> ) | The name of the file to be written.                                     |                  |  |
| <b>Input</b><br>( <i>Input</i> )        | The input filter/source whose output dataset is to written to the file. |                  | The selected object must be the result of the following: sources (includes readers), filters.<br><br>The selected dataset must be one of the following types (or a subclass of one of them): vtkDataSet. |

# How to build/compile/install

---

## Introduction

This page describes how to build and install ParaView. It covers both the release and the development versions. It covers both Unix-type systems (Linux, HP-UX, Solaris, Mac) as well as Windows.

## Prerequisites

- The ParaView build process requires CMake <sup>[1]</sup> version 2.8.2 or higher and a working compiler. On Unix-like operating systems, it also requires Make, while on Windows it requires Visual Studio (8 or later).
- Building ParaView's user interface requires TrollTech's Qt, version 4.6.\* (4.6.2 recommended) Qt is dual licensed. To compile ParaView, either the open or commercial version may be used. If you intend to make changes the ParaView's GUI and distribute the results, you must obtain a commercial license. The open source version of Qt can be found here [2]. Commercial licenses can be purchased directly from TrollTech [3]. For more information on what can be done with the open source version, read this [4].
- In order to run ParaView in parallel, MPI [5], [6] is also required.
- In order to use scripting, python is required [7].

## Download And Install CMake

CMake is a tool that makes cross-platform building simple. On several systems it will probably be already installed. If it is not, please use the following instructions to install it. If CMake does not exist on the system, and there are no pre-compiled binaries, use the instructions below on how to build it. Use the most recent source or binary version of CMake from the CMake web site.

## Using Binaries

There are several precompiled binaries available at the CMake download page <sup>[8]</sup>.

---

**On Unix-like operating systems**

Let's say on Linux, download the appropriate version and follow these instructions:

- Download [9]

```
cd $HOME
wget http://www.cmake.org/files/v2.8/cmake-2.8.2-Linux-i386.tar.gz
mkdir software
cd software
tar xvfz ../cmake-2.8.2-Linux-i386.tar.gz
```

- Now you have the directory **\$HOME/software/cmake-2.8.2-Linux-i386/bin**, and inside there are executables **cmake** and **ccmake**.
- You can also install CMake in the **/usr/local** or **/opt** by untaring and copying sub-directories. The rest of the instructions will assume the executables are in your **\$PATH**.

**On Windows**

- Download the installer [10]
- Follow the installation instructions

**On Windows, if you are not administrator**

- Download [11]
- Uncompress into some directory
- Optional: create a shortcut on the desktop.

**Build Your Own CMake****On Unix-like operating systems**

Download the source code [12]

```
cd $HOME
wget http://www.cmake.org/files/v2.8/cmake-2.8.2.tar.gz
tar xvfz cmake-2.8.2.tar.gz
cd cmake-2.8.2
./configure --prefix=$HOME/software
make
make install
```

- Again, you can install it in **/usr/local** or **/opt** by changing the prefix.

**On Windows**

To build CMake on windows, a previous version of CMake is required. This can be downloaded from the CMake download page [8].

**Download And Install Qt**

ParaView uses Trolltech's Qt as its GUI library. Qt is required whenever the ParaView client is built. See this page for Qt licensing information: [4].

- As stated above, The open source version of Qt can be found at [2].
  - For sourcecode, use the latest stable version of qt-everywhere-opensource-src-VERSION.[tar.gz or zip or dmg]. If this gives you trouble, version 4.6.2 is known to work.
  - For binaries, use the latest stable version of qt-PLATFORM-opensource-VERSION.[tar.gz or zip or dmg]. If this gives you trouble, version 4.6.2 is known to work.
- Commercial licenses can be purchased directly from TrollTech [3].



## Download And Install ffmpeg (.avi) movie libraries

When the ability to write .avi files is desired, and writing these files is not supported by the OS, ParaView can attach to an ffmpeg library. This is generally true for Linux. Ffmpeg library source code is found here: [13]

## Download And Install MESA 3D libraries

ParaView uses the OpenGL graphics drivers and card from a user's workstation. When you want to run ParaView's servers on a platform that does not include hardware OpenGL support, you must use MESA to emulate this hardware in software. Mesa is open source, and it can be downloaded from here: [14].

There is a known problem with MESA version 7.8.2 and ParaView. This has been reported to the MESA team. Version 7.7.1 has been tested and seems to work correctly as well as 7.9.

- Build as follows:
- make realclean
- make TARGET (for instance, make linux-x86-64)

Note - some platforms will complain during ParaView compiles about needing fPIC. In the configs directory, copy your platform file to another custom file, edit it, and add -fPIC to the compile lines. For instance, cp linux-x86-64 linux-x86-64-fPIC.

## Download ParaView Source Code

If you are trying to build a ParaView release, download it from the release page. For the development version, please follow the instructions below for checking it out from git.

## Download The Release

Don't forget that you can always just download the binaries from the ParaView download page <sup>[15]</sup>. This page contains binaries for several platforms and the source code for the releases.

### Note: debian build

List of packages to build ParaView on Debian:

libphonon-dev libphonon4 libqt4-assistant qt4-dev-tools libqt4-core libqt4-gui qt4-qmake libxt-dev g++ gcc  
cmake-curses-gui libqt4-opengl-dev mesa-common-dev

With MPI (using openmpi, you can use any other flavour):

openmpi-common openmpi-bin libopenmpi-dev

With Python:

python2.5-dev

## Checkout Development Version from git

Note that you may need to download and install a git client, here: [16]

### On Unix-like operating systems

```
mkdir $HOME/projects
cd $HOME/projects

cd 3.10.0
git clone git://paraview.org/ParaView.git ParaView
cd ParaView
git checkout -b trunk origin/master
git submodule init
git submodule update

Update
git fetch origin
git rebase origin/master
git submodule update
```

### On Windows

We recommend [msysgit[17]]. Msysgit provides an msys shell that has the appropriate environment set up for using git and it's tools.

## Configure ParaView With CMake

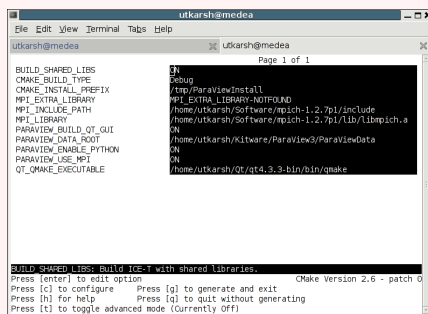
- Always use a **separate build directory**. Do not build in the source directory.

### On Unix-like systems

- Use cmake (Curses CMake GUI) from the CMake installed location. CCMake is a Curses based GUI for CMake. To run it go to the build directory and specify as an argument the src directory.

```
mkdir $HOME/projects/ParaView-bin
cd $HOME/projects/ParaView-bin
```

```
cmake $HOME/projects/ParaView3
```

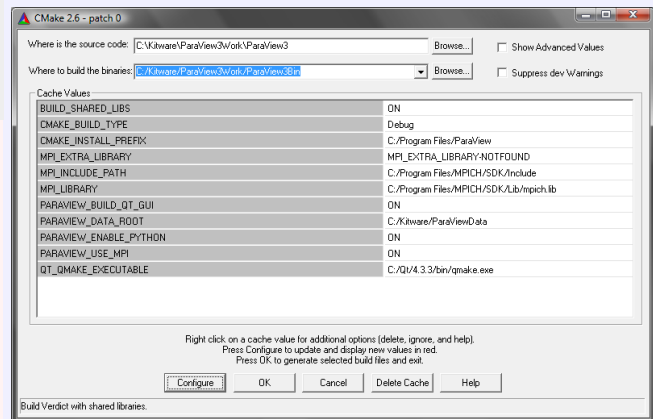


### About CCMake (Curses CMake GUI)

- Iterative process
  - Select values, run configure (c key)
  - Set the settings, run configure, set the settings, run configure, etc.
- Repeat until all values are set and the generate option is available (g key)
- Some variables (advanced variables) are not visible right away
- To see advanced variables, toggle to advanced mode (t key)
- To set a variable, move the cursor to the variable and press enter
  - If it is a boolean (ON/OFF) it will flip the value
  - If it is string or file, it will allow editing of the string
  - For file and directories, the <tab> key can be used to complete
- To search for a variable press '/' key, and to repeat the search, press the 'n' key.

### On Windows

- Use CMakeSetup from the CMake install location.
- Make sure to select the appropriate source and the build directory
- Also, make sure to pick the appropriate generator (on Visual Studio 6, pick the *Visual Studio 6* generator). Some CMake versions will ask you to select the generator the first time you press Configure instead of having a drop-down menu in the main dialog.



### About CMakeSetup (Windows CMake GUI)

- Iterative process
  - Select values, press the Configure button
  - Set the settings, run configure, set the settings, run configure, etc.
- Repeat until all values are set and the OK button becomes available.
- Some variables (advanced variables) are not visible right away
- To see advanced variables, toggle to advanced mode ("Show Advanced Values" toggle).
- To set the value of a variable, click on that value.
  - If it is boolean (ON/OFF), a drop-down menu will appear for changing the value.
  - If it is file or directory, an ellipsis button will appear ("...") on the far right of the entry. Clicking this button will bring up the file or directory selection dialog.
  - If it is a string, it will become an editable string.

## ParaView Settings

| Variable               | Description  |
|------------------------|--|
| BUILD_SHARED_LIBS      | If ON, use shared libraries. This way executables are smaller, but you have to make sure the shared libraries are on every system on the cluster.  |
| PARAVIEW_USE_MPI       | Turn this to ON to enable MPI. Other MPI options will not be available until you turn this on.   |
| MPI_LIBRARY            | Path to the MPI library (such as /usr/lib/libmpi.so). Should be found by default, but you may have to set it. (see the note below)   |
| MPI_EXTRA_LIBRARY      | Path to extra MPI library (such as /usr/lib/libmpi++.so). If the MPI distribution is MPICH, this may not be found; in this case, it is ok for this variable to be set to MPI_EXTRA_LIBRARY-NOTFOUND.   |
| MPI_INCLUDE_PATH       | Path to MPI includes (such as /usr/include/mpi). Again, this should be found by default.   |
| PARAVIEW_ENABLE_PYTHON | Makes python client scripting and the python programmable filter available.  |
| PARAVIEW_BUILD_QT_GUI  | Flag to enable/disable the building of the ParaView Qt-based client. This option is useful when building paraview on server nodes or when we are only interested in the python client, as it avoids building of the Qt client thus does not require Qt. ON by Default. |
| QT_QMAKE_EXECUTABLE    | Path to Qt's qmake executable (such as /usr/local/Trolltech/Qt-4.2.2/bin/qmake). Cmake uses this to locate the rest of the required Qt executables, headers and libraries.   |
| VTK_USE_CARBON         | For Mac, this is the default. Neither X11 nor COCOA frameworks are supported.  |
| VTK_OPENGL_HAS_OSMESA  | Turn this to ON to enable MESA (i.e., software rendering). All mesa directory and file settings below depend on this flag.   |
| OSMESA_INCLUDE_DIR     | Set this to the include directory for MESA.  |
| OPENGL_INCLUDE_DIR     | Set this to the include directory for MESA.  |
| OPENGL_gl_LIBRARY      | Set this to the libGL.a or libGL.so file.  |
| OPENGL_glu_LIBRARY     | Set this to the libGLU.a or libGLU.so file.  |
| OSMESA_LIBRARY         | Set this to the libOSMesa.a or libOSMesa.so file.  |

**Note for MPI settings:** If your MPI variables aren't set automatically (usually the case if the compiler wrapper [mpicxx] is not in the path or in some standard directory), toggle advanced options and set MPI\_COMPILER variable to the full path of your mpi compiler (usually mpicxx), and configure. This should set all the required MPI variables. If not, then you might need to enter them manually.

## Finish Configuring ParaView

### Using CMake

- Once all configuration options are set, you should be able to just run <generate> (g key).

### Using CMakeSetup

- Once all configuration options are set, you should be able to just run <generate>, by clicking the "OK" button.

## Build ParaView

You can now build ParaView using the appropriate build system.

## Using Make

CMake will now generate Make files. These makefiles have all dependencies and all rules to build ParaView on this system. You should not however try to move the build directory to another location on this system or to another system.

Once you have makefiles you should be able to just type:

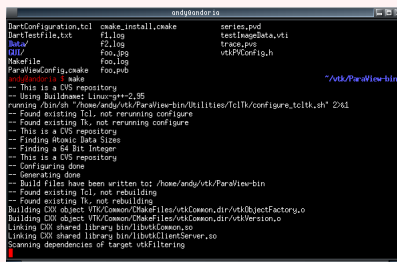
```
make
```

- If you are on multi-processor system (let's say 4 processor), you can type:

```
make -j 4
```

- To create online documentation, type the following. (Note that ParaView build does not create this documentation by default).

```
make HTMLDocumentation
```



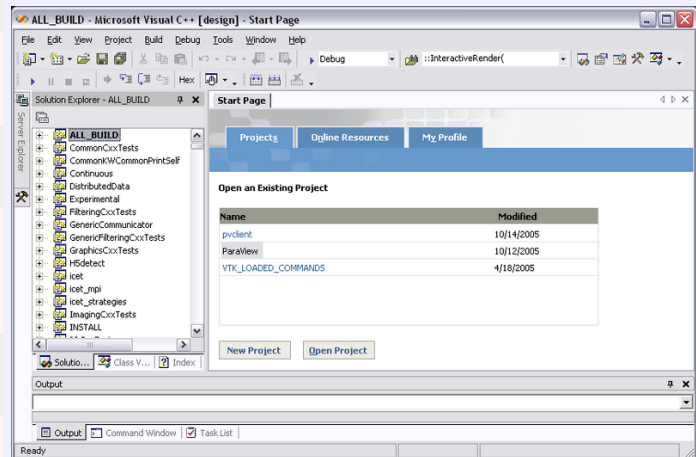
```

$ make
CMakeFiles/Makefile2:104: recipe for target 'all' failed
make: *** [all] Error 2

```

## Using Visual Studio

CMake will now create Visual Studio project files. Before you open Visual Studio, be sure that the Qt .dlls are in your path. You should now be able to open the **ParaView** project (or workspace) file. Make sure to select the appropriate build type (Debug, Release, ...). To build ParaView, simply build the **ALL\_BUILD** target. The ParaView build does not create the online documentation by default. Select the **HTMLDocumentation** target, right click, and select Build.



## Install ParaView

ParaView can be run directly from the build directory. That said, for production environments, it should be installed in some system location. For that purpose simply follow the instructions for "Distributing ParaView" and the unpack the generated archive at the appropriate install location.

## Distributing ParaView

It is possible to create distribution binaries for ParaView using **CPack** which is included in CMake. The packaged binary can be in any of the variable supported generator formats eg. DEB (debian packages), RPM (RPM Packages), NSIS (Null Soft Installer) etc. Please refer to [[18]] for a complete list.

## CMake Variables

When building ParaView so that it can be distributed, in most typical cases, ensure that the values for the following CMake variables are set correctly.

| Variable                              | Value   | Description  |
|---------------------------------------|---------|--|
| BUILD_SHARED_LIBS                     | ON      | Enables shared libraries (unless you are sure you want to do static builds)  |
| CMAKE_BUILD_TYPE                      | Release | Unless you want to end up with debug install, set this to Release  |
| PARAVIEW_USE_MPI                      | OFF     | Unless you are taking extra steps to ensure that the clients have the required MPI libraries, set this to OFF, since ParaView does not include rules to package MPI. |
| PARAVIEW_ENABLE_PYTHON                | OFF     | Unless you are taking extra steps to ensure that clients have the required python libraries/modules.   |
| PARAVIEW_GENERATE_PROXY_DOCUMENTATION | ON      | To ensure that the html documentation for the sources/filters/readers/writes is generated and packed into the application, turn the following flag ON.               |
| PARAVIEW_INSTALL_DEVELOPMENT          | OFF/ON  | To package development files so that people can build plugins/custom-apps using the installation.  |
| PARAVIEW_BUILD_QT_GUI                 | ON/OFF  | To enable building the Qt client   |

**PARAVIEW\_EXTRA\_INSTALL\_RULES\_FILE: Adding custom install rules**

You can add additional install rules by writing a custom cmake file with install rules and then set PARAVIEW\_EXTRA\_INSTALL\_RULES\_FILE cmake variable to point to that file. ParaView process all install rules that have the COMPONENT set to either "Runtime" (for executables), or "RuntimeLibraries" (for shared libraries), or "BrandedRuntime" (for custom-application specific libraries/executables) only (and "Development" when PARAVIEW\_INSTALL\_DEVELOPMENT is TRUE). So you want to add your install rules with appropriate COMPONENT values in this custom cmake file.

e.g. you want to include mpilibs in the installation, in that case the custom install rules file could look something as follows:

```
<font color="#000"><font color="#000000">
  INSTALL (FILES ${MPI_LIBRARY} ${MPI_EXTRA_LIBRARY}
           DESTINATION ${PV_INSTALL_LIB_DIR}
           COMPONENT RuntimeLibraries)
</font></font>
```

**Generate Package**

Following the configuration, simply run 'make' to compile and build.

|   |   |
|---|---|
| On Unix-like operating systems  | On Windows  |
| To generate a tar ball, in the binary directory, simply run   | To generate a null-soft installer (requires [19]), in the binary directory, simply run  |
| <pre>cpack -G TGZ --config {ParaViewBuild}/Applications/ParaView/CPackParaViewConfig.cmake</pre>                                    | <pre>cpack -G NSIS -C Release --config {ParaViewBuild}/Applications/ParaView/CPackParaViewConfig.cmake</pre>                        |
| If PARAVIEW_BUILD_QT_GUI was OFF, then to generate a package consisting of the server executables alone, use the following command: | If PARAVIEW_BUILD_QT_GUI was OFF, then to generate a package consisting of the server executables alone, use the following command: |
| <pre>cpack -G TGZ --config {ParaViewBuild}/Applications/ParaView/CPackParaViewServersConfig.cmake</pre>                             | <pre>cpack -G NSIS -C Release --config {ParaViewBuild}/Applications/ParaView/CPackParaViewServersConfig.cmake</pre>                 |

On Mac

To generate an app bundle that can be distributed, use

```
cpack -G DragNDrop --config  
{ParaViewBuild}/Applications/ParaView/CPackParaViewConfig.cmake
```

If PARAVIEW\_BUILD\_QT\_GUI was OFF, then to generate a tarball consisting of the server executables alone, use the following command:

```
cpack -G TGZ --config  
{ParaViewBuild}/Applications/ParaView/CPackParaViewServersConfig.cmake
```

## Miscellaneous Comments

- Don't untar the packaged tar ball over the build directory itself. You can always remove the build directory and untar the tar ball in its place.
- Don't use **make install** as it may install more files than required.
- CMAKE\_INSTALL\_PREFIX is not applicable when using cpack and will have no effect.
- Build trees of ParaView on non-Windows systems, always have RPATH information embedded in the binaries. When a make install is performed or CPACK is used, all RPATH information is stripped from the binaries in the install tree. By default ParaView builds forwarding executables (launchers) that are installed in the bin directory. These binaries properly setup the environment to launch the equivalent executable in the lib/paraview-x.y directory.
- If you are compiling a MESA version of the ParaView server, start the server with the --use-offscreen-memory flag.

## Notes

### Compiling on the Mac

To compile on the Mac, follow the instructions for Unix. The recommended configuration settings are: BUILD\_SHARED\_LIBS=ON.

### Environment Variables

If you build with shared libraries, you may have to add the Qt directory to your PATH environment variables to run paraview. On windows one way to do so is to open up the environment variables dialog by clicking through Start->Control Panel->System->Advanced->Environment Variables. From that dialog add a new user variable called PATH with a value of C:\Qt\4.6.3\bin. For other operating systems, add Qt/4.6.3/lib to your LD\_LIBRARY\_PATH environment variable.

---

## References

- [1] <http://www.cmake.org>
- [2] <ftp://ftp.trolltech.no/qt/source/>
- [3] <http://trolltech.com>
- [4] [http://www.paraview.org/Wiki/ParaView\\_III\\_and\\_Qt\\_licensing](http://www.paraview.org/Wiki/ParaView_III_and_Qt_licensing)
- [5] <http://www-unix.mcs.anl.gov/mpi/>
- [6] <http://www.lam-mpi.org/>
- [7] <http://www.python.org>
- [8] <http://www.cmake.org/HTML/Download.html>
- [9] <http://www.cmake.org/files/v2.8/cmake-2.8.2-Linux-i386.tar.gz>
- [10] <http://www.cmake.org/files/v2.8/cmake-2.8.2-win32-x86.exe>
- [11] <http://www.cmake.org/files/v2.8/cmake-2.8.2-win32-x86.zip>
- [12] <http://www.cmake.org/files/v2.8/cmake-2.8.2.tar.gz>
- [13] <http://www.ffmpeg.org/>
- [14] <http://www.mesa3d.org/>
- [15] <http://www.paraview.org/HTML/Download.html>
- [16] <http://git-scm.com/>
- [17] <http://code.google.com/p/msysgit>
- [18] <http://www.cmake.org/Wiki/CMake:CPackPackageGenerators>
- [19] [http://nsis.sourceforge.net/Main\\_Page](http://nsis.sourceforge.net/Main_Page)

---

# How to write parallel VTK readers

## Writing ParaView Readers

If the format of your data files is not one supported by default in ParaView (see section Error: Reference source not found), you will either need to convert your files to a format ParaView can read, or you must write your own data file reader for ParaView. The reader must operate within a standard VTK pipeline. In this chapter, we will discuss integrating the new reader class into VTK, including outlining which C++ methods should be implemented for the reader to work properly. The necessary user interface and server manager XML will be described. Creating parallel readers and readers that output multiple parts will also be covered. For VTK information beyond the scope of the chapter, see *The VTK User's Guide* by Kitware, Inc.

## Integrating with VTK

VTK is written in C++, and new readers should also be written in this language. A reader plays the role of a source in a VTK pipeline, and must be implemented as a class deriving from `vtkAlgorithm` or one of its subclasses. The best choice for the immediate superclass of a new reader depends on the reader's output type. For example, a reader producing an instance of `vtkPolyData` may derive from `vtkPolyDataAlgorithm` to simplify its implementation. In order for a reader to function properly within VTK's pipeline mechanism, it must be able to respond to standard requests. This is implemented by overriding one or more of the following methods from the chosen superclass.

The interface to these methods utilizes `vtkInformation` objects which are heterogeneous maps storing key/value pairs. Many of these methods have the same three arguments. The first argument is of type `vtkInformation*` and contains at least one key specifying the request itself. The second argument is of type `vtkInformationVector**` and stores information about the input connections to the algorithm. This can be ignored by readers because they have no input connections. The third argument is of type `vtkInformationVector*` and contains one `vtkInformation` object for each

---



output port of the algorithm. Most readers will have only one output port, but some may have multiple output ports (see the next section). All output information and data from the reader will be stored in one of these information objects.

**ProcessRequest:** This method is the entry point into a `vtkAlgorithm` through which the pipeline makes requests. A reader may override this method and implement responses to all requests. The method should be placed in the public section of the reader class. It should return 1 for success and 0 for failure. Full documentation of this method is beyond the scope of this chapter. Most readers should derive from one of the output-type-specific classes and implement the request-specific methods described below.

**RequestInformation:** This method is invoked by the superclass's **ProcessRequest** implementation when it receives a `REQUEST_INFORMATION` request. In the output port, it should store information about the data in the input file. For example, if the reader produces structured data, then the whole extent should be set here (shown below).

```
int vtkExampleReader::RequestInformation(
    vtkInformation*, vtkInformationVector**,
    vtkInformationVector* outVec)
{
    vtkInformation* outInfo = outVec->GetInformationObject(0);
    int extent[6];

    // ... read file to find available extent ...

    //store that in the pipeline
    outInfo->Set
        (vtkStreamingDemandDrivenPipeline::WHOLE_EXTENT(), extent, 6);

    // ... store other information ...

    return 1;
}
```

This method is necessary when configuring a reader to operate in parallel. (This will be further discussed later in this chapter.) It should be placed in the protected section of the reader class. The method should return 1 for success and 0 for failure.

**RequestData:** This method is invoked by the superclass's **ProcessRequest** implementation when it receives a `REQUEST_DATA` request. It should read data from the file and store it in the corresponding data object in the output port. The output data object will have already been created by the pipeline before this request is made. The amount of data to read may be specified by keys in the output port information. For example, if the reader produces `vtkImageData`, this method might look like this.

```
int vtkExampleReader::RequestData(
    vtkInformation*, vtkInformationVector**,
    vtkInformationVector* outVec)
{
    vtkInformation* outInfo = outVec->GetInformationObject(0);
    vtkImageData* outData = vtkImageData::SafeDownCast
        (outInfo->Get(vtkDataObject::DATA_OBJECT()));

    int extent[6] = {0,-1,0,-1,0,-1};
```

```

outInfo->Get
    (vtkStreamingDemandDrivenPipeline::UPDATE_EXTENT(), extent);

outData->SetExtent(extent);

// ... read data for this extent from the file ...

return 1;
}

```

The method should be placed in the protected section of the reader class. It should return 1 for success and 0 for failure.

**CanReadFile:** The purpose of this method is to determine whether this reader can read a specified data file. Its input parameter is a `const char*` specifying the name of a data file. In this method you should not actually read the data but determine whether it is the correct format to be read by this reader. This method should return an integer value: 1 indicates that the specified file is of the correct type; 0 indicates it is not. It is not absolutely required that this method be implemented, but ParaView will make use of it if it exists.

**SetFileName:** This method allows you to specify the name of the data file to be loaded by your reader. The method is not required to have this exact name, but a method with this functionality must be implemented. The easiest way to implement `SetFileName` is with a `vtkSetStringMacro` in the header file for this class. (There is also an associated `vtkGetStringMacro` for implementing `GetFileName`.) This method handles allocating an array to contain the file name and lets the reader know that the pipeline should be updated when the name is changed.

```

vtkSetStringMacro (FileName);

```

When using this macro, you must also add a `FileName` instance variable of type `char*` in the protected section of this class. In the constructor for your reader, assign `FileName` the value `NULL` before you use `SetFileName` for the first time. In the destructor for your reader, call `SetFileName(0)` to free the file name storage.

## Multi-Group (Multi-Block and AMR) Readers

As of VTK 5.0 and ParaView 2.4, multi-block and AMR datasets are supported. Multi-group readers follow the same guidelines as described in the previous sections. For convenience, you can subclass multi-block readers from `vtkMultiBlockDataSetAlgorithm` and AMR readers from `vtkHierarchicalDataSetAlgorithm`. If you do not subclass from one of these classes, make sure to implement the `CreateDefaultExecutive()`, `FillOutputPortInformation()`, and `FillInputPortInformation()` methods appropriately. (You can use `vtkMultiBlockDataSetAlgorithm` as a starting point.) Two good examples of multi-group dataset readers are `vtkMultiBlockPLOT3DReader` and `vtkXMLHierarchicalDataReader`.

## Parallel Readers

Unless otherwise specified, a VTK reader used in ParaView will cause the entire data set to be read on the first process. ParaView will then redistribute the data to the other processes. It is more desirable to have ParaView do the reading in parallel as well so that the data set is already appropriately divided across the processors. Changes should be made in two methods for ParaView readers to operate in parallel: `RequestInformation` and `RequestData`. Exactly which changes should be made depends on whether structured or unstructured data set types are to be read.

## Structured

In `RequestInformation` for readers of structured data (`vtkStructuredGrid`, `vtkRectilinearGrid`, or `vtkImageData`), the `WHOLE_EXTENT` key should be set on the output information object, so that downstream filters can take into account the overall dimensions of the data. The `WHOLE_EXTENT` is specified using six parameters: the minimum and maximum index along each of the three coordinate axes ( $i_{\min}$ ,  $i_{\max}$ ,  $j_{\min}$ ,  $j_{\max}$ ,  $k_{\min}$ ,  $k_{\max}$  or a single int array of length 6) for the entire data set. Example C++ code demonstrating this is shown below.

```
int vtkDEMReader::RequestInformation (
    vtkInformation * vtkNotUsed(request),
    vtkInformationVector ** vtkNotUsed( inputVector ),
    vtkInformationVector *outputVector)
{
    vtkInformation* outInfo =

    outputVector->GetInformationObject(0);
    int extent[6];

    //Read entire extent from file.
    ...

    outInfo->Set
        (vtkStreamingDemandDrivenPipeline::WHOLE_EXTENT(), extent, 6);

    return 1;
}
```

Before doing any processing in the `RequestData` method, first get the `UPDATE_EXTENT` from the output information object. This key contains the sub-extent of the `WHOLE_EXTENT` for which the current process is responsible. The current process is responsible for filling in the data values for the update extent that is returned. An example of doing this is shown below.

```
int vtkDEMReader::RequestData (
    vtkInformation* vtkNotUsed( request ),
    vtkInformationVector** vtkNotUsed( inputVector ),
    vtkInformationVector* outputVector)
{
    // get the data object and find out what part we need to
    // read now
    vtkInformation *outInfo = outputVector->GetInformationObject(0);
    int subext[6];
    outInfo->Get
        (vtkStreamingDemandDrivenPipeline::UPDATE_EXTENT(),
        subext);

    //read that part of the data in from the file
    //and put it in the output data
    ...

    return 1;
}
```

```
}
```

## Unstructured

In the unstructured case (`vtkPolyData` or `vtkUnstructuredGrid`), the `MAXIMUM_NUMBER_OF_PIECES` key should be set on the output information object in `RequestInformation`. This specifies the maximum number of pieces that can be generated from the file. If this reader can only read the entire data set, then the maximum number of pieces should be set to 1. If the input file can be read into as many pieces as needed (i.e., one per processor), the maximum number of pieces should be set to -1, as shown below.

```
outInfo->Set(vtkStreamingDemandDrivenPipeline::MAXIMUM_NUMBER_OF_PIECES(),
-1);
```

In the `RequestData` method, the reader should first get the `UPDATE_NUMBER_OF_PIECES` and `UPDATE_PIECE_NUMBER` from the output information object. The value returned from getting the `UPDATE_NUMBER_OF_PIECES` specifies the number of pieces into which the output data set will be broken. Getting `UPDATE_PIECE_NUMBER` returns the piece number (0 to `UPDATE_NUMBER_OF_PIECES`-1) for which the current process is responsible. The reader should use this information to determine which part of the data set the current process should read. Example code that demonstrates this is shown below.

```
int vtkUnstructuredGridReader::RequestData(
    vtkInformation *,
    vtkInformationVector **,
    vtkInformationVector *outputVector)
{
    vtkInformation *outInfo = outputVector->GetInformationObject(0);
    int piece, numPieces;

    piece = outInfo->Get
        (vtkStreamingDemandDrivenPipeline::UPDATE_PIECE_NUMBER());

    numPieces = outInfo->Get
        (vtkStreamingDemandDrivenPipeline::UPDATE_NUMBER_OF_PIECES());

    //skip to proper offset in the file and read piece
    ...

    return 1;
}
```

It is possible that your data file can only be broken into a specified number of pieces, and that this number is different than the number of processors being used (i.e., the result of getting `UPDATE_NUMBER_OF_PIECES`). If the number of processors is larger than the possible number of pieces, then each processor beyond the number of available pieces should produce an empty output by calling `Initialize()` on the output. If the number of processors is smaller than the number of pieces, you should internally redistribute the extra data across the processors. For example, if your data set can produce ten pieces, and you are using five processors for reading, then process 0 could read pieces 0 and 5; process 1 could read pieces 1 and 6; etc.

## Required XML

To use your new reader within ParaView, you must write XML code for the server manager and for the client. The server-side XML for ParaView's readers is located in the file `Servers/ServerManager/Resources/readers.xml`. When ParaView is compiled, the server manager XML code for a reader is used to create a proxy object for it. The ParaView client accesses the VTK class for the reader on the server through the proxy. Below is an excerpt from the server manager XML code for `vtkXMLPolyDataReader`. A few parts of this XML code segment are particularly worth noting, and more extensive information about server manager XML code can be found in section [Error: Reference source not found](#).

First, notice the `StringVectorProperty` element named `"FileName"`. It has a command attribute called `"SetFileName"`. ParaView uses this property to tell the reader what file it should examine. This is done very early in the lifetime of a reader. Typically, ParaView code will give the reader a filename and then call the reader's `CanReadFile` method. If `CanReadFile` succeeds, ParaView will first call `RequestInformation` to get general information about the data within the file, and then call `RequestData` to read the actual data and produce a `vtkDataObject`.

The second notable portion of the XML code below is two properties that let the user choose particular arrays to load from the data file. When ParaView sees them, it creates a section on the **Properties** tab for the reader that lets the user select from the available cell-centered data arrays. In the following discussion, one can simply replace `'cell'` with `'point'` in order to let the user choose from the available point-centered arrays.

The `StringVectorProperty` named `"CellArrayStatus"` lets the ParaView client call the `SetCellArrayStatus` method on the server. The `SetCellArrayStatus` method is how the reader is told what it should do with each of the arrays the data file contains. It takes two parameters; the first is the name of the array, and the second is an integer indicating whether to read the array (1) or not (0).

The `StringVectorProperty` named `'CellArrayInfo'` is an information property; that is, one through which the ParaView client gathers information from the server. The ParaView client uses it to gather the names of the cell-centered arrays from the reader. (See the nested `Property` sub-element in `CellArrayStatus's` `ArraySelectionDomain` element.) In order for this collection to work, two methods be implemented in the reader. `GetNumberOfCellArrays` returns the number (an int) of cell-centered arrays in the data file. This method does not accept any parameters. `GetCellArrayName` takes a single parameter the array's index (starting from 0). This method returns the name of the array (a `const char*`) with this index or `NULL` if the array has no name or if the index is larger than the number of arrays.

```
<SourceProxy name="XMLPolyDataReader"
              class="vtkXMLPolyDataReader"
              label="XML Polydata reader">
  <StringVectorProperty name="FileName"
    command="SetFileName"
    animateable="0"
    number_of_elements="1">
    <FileListDomain name="files"/>
  </StringVectorProperty>

  <StringVectorProperty name="CellArrayInfo"
    information_only="1">
    <ArraySelectionInformationHelper attribute_name="Cell"/>
  </StringVectorProperty>

  <StringVectorProperty name="CellArrayStatus"
    command="SetCellArrayStatus"
    number_of_elements="0"
```

```

repeat_command="1" number_of_elements_per_command="2"
element_types="2 0"
information_property="CellArrayInfo"
label="Cell Arrays">
  <ArraySelectionDomain name="array_list">
    <RequiredProperties>
      <Property name="CellArrayInfo"
        function="ArrayList"/>
    </RequiredProperties>
  </ArraySelectionDomain>
</StringVectorProperty>

</SourceProxy>

```

The client-side XML is extremely simple. The purpose of the client-side XML is to enable readers from the server-side XML and to associate file extensions with them. In the case where multiple readers can read files that have the same file extensions, the `CanReadFile` method is called on each one in order to choose the correct one for each data file. The client-side XML file for readers is located in `Qt/Components/Resources/XML/ParaViewReaders.xml`. The portion of that file related to `XMLPolyDataReader` follows.

```

<Reader name="XMLPolyDataReader"
  extensions="vtp"
  file_description="VTK PolyData Files">
</Reader>

```

An alternative to modifying the ParaView source code directly is to use ParaView's plugin architecture. With a plugin, the same C++ source code and XML content must be written, but it is kept outside of the ParaView source code proper and is compiled separately from ParaView. Plugins are discussed in chapter Error: Reference source not found.

# Article Sources and Contributors

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